

	Final Report
	Re-Issued Repor
V	Revised Report
,	port Date: Apr-17 16:19

Laboratory Report

Gulf Oil L.P. 281 Eastern Avenue Chelsea, MA 02150 Attn: Andrew P. Adams

Project: Gulf Terminal - Chelsea, MA

Project #: Gulf Chelsea

Laboratory ID	Client Sample ID	<u>Matrix</u>	Date Sampled	Date Received
SC32731-01	Chelsea Creek	Surface Water	23-Mar-17 10:00	23-Mar-17 15:00
SC32732-01	Outfall 003	Surface Water	23-Mar-17 10:00	23-Mar-17 15:00

I attest that the information contained within the report has been reviewed for accuracy and checked against the quality control requirements for each method. These results relate only to the sample(s) as received.

All applicable NELAC requirements have been met.

Massachusetts # M-MA138/MA1110 Connecticut # PH-0777 Florida # E87936 Maine # MA138 New Hampshire # 2972/2538 New Jersey # MA011 New York # 11393 Pennsylvania # 68-04426/68-02924 Rhode Island # LAO00348 USDA # P330-15-00375 Vermont # VT-11393



Authorized by:

June O'Connor Laboratory Director

Eurofins Spectrum Analytical holds primary certification in the State of Massachusetts for the analytes as indicated with an X in the "Cert." column within this report. Please note that the State of Massachusetts does not offer certification for all analytes. Please refer to our website for specific certification holdings in each state.

Please note that this report contains 23 pages of analytical data plus Chain of Custody document(s). When the Laboratory Report is indicated as revised, this report supersedes any previously dated reports for the laboratory ID(s) referenced above. Where this report identifies subcontracted analyses, copies of the subcontractor's test report are available upon request. This report may not be reproduced, except in full, without written approval from Eurofins Spectrum Analytical, Inc.

Eurofins Spectrum Analytical, Inc. is a NELAC accredited laboratory organization and meets NELAC testing standards. Use of the NELAC logo however does not insure that Eurofins Spectrum Analytical, Inc. is currently accredited for the specific method or analyte indicated. Please refer to our Quality web page at www.spectrum-analytical.com for a full listing of our current certifications and fields of accreditation. States in which Eurofins Spectrum Analytical, Inc. holds NELAC certification are New York, New Hampshire, New Jersey, Pennsylvania and Florida. All analytical work for Volatile Organic and Air analysis are transferred to and conducted at our 830 Silver Street location (PA-68-04426).

Please contact the Laboratory or Technical Director at 800-789-9115 with any questions regarding the data contained in this laboratory report.

CASE NARRATIVE:

Data has been reported to the MDL. This report includes estimated concentrations detected below the RDL and above the MDL (J-Flag).

All non-detects and all results below the detection limit are reported as "<" (less than) the detection limit in this report.

The samples were received 2.9 degrees Celsius, please refer to the Chain of Custody for details specific to temperature upon receipt. An infrared thermometer with a tolerance of +/- 1.0 degrees Celsius was used immediately upon receipt of the samples.

If a Matrix Spike (MS), Matrix Spike Duplicate (MSD) or Duplicate (DUP) was not requested on the Chain of Custody, method criteria may have been fulfilled with a source sample not of this Sample Delivery Group.

Analyses for Total Hardness, pH, and Total Residual Chlorine fall under the state of Pennsylvania code Chapter 252.6 accreditation by rule.

Please note this work order contains 24 pages of analytical data from New England Bioassay.

Report Re-issue Case Narrative 4/11/17:

The laboratory report for SC32731/32 was re-issued to report Phenol by 8270D only.

See below for any non-conformances and issues relating to quality control samples and/or sample analysis/matrix.

EPA 200.7

Duplicates:

1705143-DUP1 Source: SC32731-01

The Reporting Limit has been raised to account for matrix interference.

Cadmium

Copper

Lead

Nickel Zinc

Samples:

SC32731-01 Chelsea Creek

The Reporting Limit has been raised to account for matrix interference.

Cadmium

Lead

Nickel

Zinc

EPA 200.8

Spikes:

1705144-PS1 Source: SC32731-01

The spike recovery was outside acceptance limits for the MS, MSD and/or PS due to matrix interference. The LCS and/or LCSD were within acceptance limits showing that the laboratory is in control and the data is acceptable.

Copper

Duplicates:

1705144-DUP1 Source: SC32731-01

EPA 200.8

Duplicates:

1705144-DUP1 Source: SC32731-01

MRL raised to correlate to batch QC reporting limits.

Copper

The Reporting Limit has been raised to account for matrix interference.

Chromium

Copper

Samples:

SC32731-01

Chelsea Creek

MRL raised to correlate to batch QC reporting limits.

Copper

The Reporting Limit has been raised to account for matrix interference.

Copper

SC32732-01

Outfall 003

The Reporting Limit has been raised to account for matrix interference.

Chromium

SM4500-Cl-G (11)

Spikes:

1705084-MS1

Source: SC32731-01

The spike recovery for this QC sample is outside the established control limits. The sample results for the QC batch were accepted based on LCS/LCSD or SRM recoveries within the control limits.

Total Residual Chlorine

1705084-MSD1

Source: SC32731-01

The spike recovery for this QC sample is outside the established control limits. The sample results for the QC batch were accepted based on LCS/LCSD or SRM recoveries within the control limits.

Total Residual Chlorine

SM4500-NH3 C. (11)

Laboratory Control Samples:

1705173 SRM

Ammonia as N percent recovery 84 (86-114) is outside individual acceptance criteria, but within overall method allowances. All reported results of the following samples are considered to have a potentially low bias:

Chelsea Creek

Outfall 003

SW846 8260C

Calibration:

1703030

Analyte quantified by quadratic equation type calibration.

Naphthalene

SW846 8260C

Calibration:

1703030

This affected the following samples:

1705104-BLK1

1705104-BS1

1705104-BSD1

Chelsea Creek

Outfall 003

S703239-ICV1

S703263-CCV1

Laboratory Control Samples:

1705104 BS/BSD

Ethanol percent recoveries (138/134) are outside individual acceptance criteria (70-130), but within overall method allowances.

All reported results of the following samples are considered to have a potentially high bias:

Outfall 003

Samples:

S703263-CCV1

Analyte percent drift is outside individual acceptance criteria (20), but within overall method allowances.

Ethanol (37.8%)

This affected the following samples:

1705104-BLK1

1705104-BS1

1705104-BSD1

Outfall 003

SW846 8270D

Samples:

S703277-CCV1

Analyte percent drift is outside individual acceptance criteria (20), but within overall method allowances.

2,4-Dinitrophenol (30.7%)

4,6-Dinitro-2-methylphenol (34.6%)

4-Nitrophenol (23.5%)

This affected the following samples:

1705073-BLK1

1705073-BS1

1705073-BSD1

SW846 8270D SIM

Laboratory Control Samples:

1705073 BSD

Indeno (1,2,3-cd) pyrene RPD 25% (20%) is outside individual acceptance criteria.

Naphthalene RPD 21% (20%) is outside individual acceptance criteria.

Samples:

SW846 8270D SIM

Samples:

S703276-CCV1

Analyte percent difference is outside individual acceptance criteria (20), but within overall method allowances.

Dibenzo (a,h) anthracene (28.1%) Indeno (1,2,3-cd) pyrene (21.4%)

This affected the following samples:

1705073-BLK2 1705073-BS2 1705073-BSD2 Chelsea Creek Outfall 003

Sample Acceptance Check Form

Client:

Gulf Oil L.P.

Project:		Gulf Terminal - Chelsea, MA / Gulf Chelsea			
Work Order		SC32731			
Sample(s) re	eceived on:	3/23/2017			
The followi	ng outlines the	condition of samples for the attached Chain of Custody upon receipt.			
			<u>Yes</u>	No	<u>N/A</u>
Wei	re custody seal	s present?		\checkmark	
Wei	re custody seal	s intact?			✓
Wei	re samples rec	eived at a temperature of $\leq 6^{\circ}$ C?	\checkmark		
Wei	re samples refi	igerated upon transfer to laboratory representative?	✓		
Wei	re sample cont	ainers received intact?	\checkmark		
		perly labeled (labels affixed to sample containers and include sample ID, site roject number and the collection date)?			
Wei	re samples acc	ompanied by a Chain of Custody document?	\checkmark		
incl	lude sample ID	stody document include proper, full, and complete documentation, which shall by, site location, and/or project number, date and time of collection, collector's name, sample matrix and any special remarks concerning the sample?			
Did	sample contain	ner labels agree with Chain of Custody document?	\checkmark		
Wei	re samples rec	eived within method-specific holding times?	\checkmark		

Sample Acceptance Check Form

Client:

Gulf Oil L.P.

Project:		Gulf Terminal - Chelsea, MA / Gulf Chelsea			
Work Or	der:	SC32732			
Sample(s	s) received on:	3/23/2017			
The follo	owing outlines the	condition of samples for the attached Chain of Custody upon receipt.			
			Yes	No	<u>N/A</u>
•	Were custody sea	s present?		\checkmark	
1	Were custody sea	s intact?			✓
•	Were samples rec	eived at a temperature of $\leq 6^{\circ}$ C?	✓		
7	Were samples ref	igerated upon transfer to laboratory representative?	✓		
,	Were sample cont	ainers received intact?	\checkmark		
		perly labeled (labels affixed to sample containers and include sample ID, site roject number and the collection date)?	\checkmark		
7	Were samples acc	ompanied by a Chain of Custody document?	✓		
j	include sample II	stody document include proper, full, and complete documentation, which shall b, site location, and/or project number, date and time of collection, collector's name, sample matrix and any special remarks concerning the sample?			
]	Did sample conta	ner labels agree with Chain of Custody document?	\checkmark		
,	Were samples rec	eived within method-specific holding times?	✓		

Summary of Hits

Lab ID: SC32731-01

Client ID: Chelsea Creek

Parameter	Result	Flag Reporting Limit	Units	Analytical Method
Copper	0.0972	R01, R(0.00800	mg/l	EPA 200.8
Salinity	28.8	1.00	ppt (1000)	SM 2520 (01)
Total Solids	32800	100	mg/l	SM2540 B (11)
Total Suspended Solids	13.5	2.5	mg/l	SM2540D (11)
Total Residual Chlorine	0.034	0.020	mg/l	SM4500-Cl-G (11)
Total Organic Carbon	2.39	1.00	mg/l	SM5310B (00, 11)

Lab ID: SC32732-01

Client ID: Outfall 003

Parameter	Result	Flag	Reporting Limit	Units	Analytical Method
Cadmium	0.0004	J	0.0025	mg/l	EPA 200.7
Copper	0.0092		0.0050	mg/l	EPA 200.7
Lead	0.0095		0.0075	mg/l	EPA 200.7
Nickel	0.0044	J	0.0050	mg/l	EPA 200.7
Zinc	0.0393		0.0050	mg/l	EPA 200.7
Chromium	0.00979	R01,	D 0.00500	mg/l	EPA 200.8
Salinity	1.70		1.00	ppt (1000)	SM 2520 (01)
Total Solids	1940		5.00	mg/l	SM2540 B (11)
Total Suspended Solids	25.0		2.5	mg/l	SM2540D (11)
Total Residual Chlorine	0.079		0.020	mg/l	SM4500-Cl-G (11)
Ammonia as N	0.420		0.200	mg/l	SM4500-NH3 C. (11)
Total Organic Carbon	5.58		1.00	mg/l	SM5310B (00, 11)

Please note that because there are no reporting limits associated with hazardous waste characterizations or micro analyses, this summary does not include hits from these analyses if included in this work order.

Sample Id Chelsea C SC32731-					Project # Chelsea	;	Matrix Surface W		ection Date -Mar-17 10			<u>ceived</u> Mar-17	
CAS No.	Analyte(s)	Result	Flag	Units	*RDL	MDL	Dilution	Method Ref.	Prepared	Analyzed	Analyst	Batch	Cert.
Volatile Or	ganic Compounds												
Volatile Or	ganic Aromatics by SW84	6 8260											
<u>Prepared</u>	by method SW846 5030 V	Vater MS											
71-43-2	Benzene	< 1.0		μg/l	1.0	0.3	1	SW846 8260C	24-Mar-17	24-Mar-17	GMA	1705104	
100-41-4	Ethylbenzene	< 1.0		μg/l	1.0	0.3	1	"	"	"	ıı		
91-20-3	Naphthalene	< 1.0		μg/l	1.0	0.4	1	"	"	"	"	"	
108-88-3	Toluene	< 1.0		μg/l	1.0	0.3	1	"	"	"	"	"	
179601-23-1	m,p-Xylene	< 2.0		μg/l	2.0	0.4	1	"	"	"	"	"	
95-47-6	o-Xylene	< 1.0		μg/l	1.0	0.3	1	"	"	"	"	"	
Surrogate r	ecoveries:												
460-00-4	4-Bromofluorobenzene	106			70-13	0 %		"		"	"	"	
2037-26-5	Toluene-d8	101			70-13			"	"	"		"	
17060-07-0	1,2-Dichloroethane-d4	106			70-13	0 %		"	"	"	"		
1868-53-7	Dibromofluoromethane	104			70-13	0 %		"	"	"	"		
Semivolati	le Organic Compounds by (GCMS											
SVOCs by	· .	Genis											
-	by method SW846 3510C												
83-32-9	Acenaphthene	< 0.051		μg/l	0.051	0.031	1	SW846 8270D SIM	24-Mar-17	25-Mar-17	MSL	1705073	
208-96-8	Acenaphthylene	< 0.051		μg/l	0.051	0.032	1	"	"	"	"		
120-12-7	Anthracene	< 0.051		μg/l	0.051	0.027	1	"	"	"	"		
56-55-3	Benzo (a) anthracene	< 0.051		μg/l	0.051	0.024	1	"	"	"	"	"	
50-32-8	Benzo (a) pyrene	< 0.051		μg/l	0.051	0.036	1		"	"	"	"	
205-99-2	Benzo (b) fluoranthene	< 0.051		μg/l	0.051	0.035	1		"	"	"	"	
191-24-2	Benzo (g,h,i) perylene	< 0.051		μg/l	0.051	0.027	1		"	"	"	"	
207-08-9	Benzo (k) fluoranthene	< 0.051		μg/l	0.051	0.028	1		"	"	"	"	
218-01-9	Chrysene	< 0.051		μg/l	0.051	0.024	1	"	"	"	"		
53-70-3	Dibenzo (a,h) anthracene	< 0.051		μg/l	0.051	0.026	1	"	"	"	"		
206-44-0	Fluoranthene	< 0.051		μg/l	0.051	0.020	1	"	"	"	"		
86-73-7	Fluorene	< 0.051		μg/l	0.051	0.030	1	"	"		"		
193-39-5	Indeno (1,2,3-cd) pyrene	< 0.051		μg/l	0.051	0.022	1	"	"		"		
91-20-3	Naphthalene	< 0.051		μg/l	0.051	0.027	1	"	"	"	"		
85-01-8	Phenanthrene	< 0.051		μg/l	0.051	0.027	1	"	"	"	"		
129-00-0	Pyrene	< 0.051		μg/l	0.051	0.022	1		"		"	"	
O													
Surrogate r		100			30-13	0.0/					"		
	Benzo (e) pyrene-d12				30-13	0%							
	ols by EPA 200/6000 Series In the Market Series In												
	Preservation	Field Preserved; pH<2 confirmed		N/A			1	EPA 200/6000 methods	24-Mar-17		BK	1705119	
	ds by EPA 200 Series Metho	ods											
7440-43-9	Cadmium	< 0.0020	U, R01, D	mg/l	0.0125	0.0020	5	EPA 200.7		31-Mar-17	TBC	1705143	Х
7440-50-8	Copper	0.0972	R01, R06, D	mg/l	0.00800	0.00089	10	EPA 200.8	"	28-Mar-17	edt	1705144	Х
7440-02-0	Nickel	< 0.0050	U, R01, D	mg/l	0.0250	0.0050	5	EPA 200.7	u	31-Mar-17	TBC	1705143	Х

Sample Id Chelsea G SC32731				Client P Gulf C			<u>Matrix</u> Surface W	· · · · · · · · · · · · · · · · · · ·	ection Date -Mar-17 10	,		<u>ceived</u> Mar-17	
CAS No.	Analyte(s)	Result	Flag	Units	*RDL	MDL	Dilution	Method Ref.	Prepared	Analyzed	Analyst	Batch	Cert.
Total Met	als by EPA 200 Series Meth	ods											
7439-92-1	Lead	< 0.0171	U, R01, D	mg/l	0.0375	0.0171	5	EPA 200.7	27-Mar-17	31-Mar-17	TBC	1705143	Х
7440-66-6	Zinc	< 0.0136	U, R01, D	mg/l	0.0250	0.0136	5	п	u	"	"	"	X
General C	Chemistry Parameters												
7782-50-5	Total Residual Chlorine	0.034		mg/l	0.020	0.006	1	SM4500-CI-G (11)	24-Mar-17 09:36	24-Mar-17 16:43	RLT	1705084	X
Prepared	by method SM4500-NH3	B (11)											
	Ammonia as N	< 0.200		mg/l	0.200	0.118	1	SM4500-NH3 C. (11)	27-Mar-17	27-Mar-17	AHK	1705173	Х
	рH	7.91		pH Units			1	ASTM D 1293-99B	23-Mar-17 17:56	23-Mar-17 18:30	BD	1705079	Х
	Salinity	28.8		ppt (1000)	1.00	0.144	1	SM 2520 (01)	28-Mar-17	28-Mar-17	BD	1705216	
	Total Solids	32,800		mg/l	100	30.6	1	SM2540 B (11)	25-Mar-17	28-Mar-17	CMB	1705113	
	Total Suspended Solids	13.5		mg/l	2.5	8.0	1	SM2540D (11)	28-Mar-17	30-Mar-17	CMB	1705207	Χ
	Total Organic Carbon	2.39		mg/l	1.00	0.246	1	SM5310B (00, 11)	30-Mar-17	30-Mar-17	RLT	1705454	X
Aquatic T Prepared	oxicity by method NA												
Analysis p	erformed by GZA Geoenviro	nmental, Inc N	Aanchester,	<i>CT*</i> -									
	Aquatic Toxicity	See Report		N/A			1	EPA-821-R-02-0 12		10-Apr-17		'[none]'	

Outfall 00				Client F	Project #		Matrix	Coll	ection Date	/Time	Re-	ceived	
					Chelsea		Surface Wa	'	-Mar-17 10			Mar-17	
SC32732-	01			Sun C			-a.1400 111						
CAS No.	Analyte(s)	Result	Flag	Units	*RDL	MDL	Dilution	Method Ref.	Prepared	Analyzed	Analyst	Batch	Cer
Volatile Or	ganic Compounds												
	ganic Compounds by SW by method SW846 5030 W												
71-43-2	Benzene	< 1.00		μg/l	1.00	0.28	1	SW846 8260C	24-Mar-17	24-Mar-17	GMA	1705104	
100-41-4	Ethylbenzene	< 1.00		μg/l	1.00	0.33	1	"	"	"	"	"	
634-04-4	Methyl tert-butyl ether	< 1.00		μg/l	1.00	0.24	1	"	"	"	"	"	
1-20-3	Naphthalene	< 1.00		μg/l	1.00	0.35	1	"	"	"	"	"	
08-88-3	Toluene	< 1.00		μg/l	1.00	0.30	1	"	"	"	"	"	
5-01-4	Vinyl chloride	< 1.00		μg/l	1.00	0.47	1	"	"	"	"	"	
79601-23-1	m,p-Xylene	< 2.00		μg/l	2.00	0.38	1	"	"	"	"	"	
5-47-6	o-Xylene	< 1.00		μg/l	1.00	0.28	1	"	"	"	"	"	
75-65-0	Tert-Butanol / butyl alcohol	< 10.0		μg/l	10.0	5.90	1	"	"	"	"	"	
4-17-5	Ethanol	< 200		μg/l	200	30.9	1	· ·	"	"	"	u u	
Surrogate r	ecoveries:												
160-00-4	4-Bromofluorobenzene	102			70-13	80 %		"	"	"	"	"	
2037-26-5	Toluene-d8	100			70-13	80 %		"	"	"	"	"	
7060-07-0	1,2-Dichloroethane-d4	102			70-13	80 %		"	"	"	"	"	
868-53-7	Dibromofluoromethane	98			70-13	80 %			"	"	"	"	
cid Extra	le Organic Compounds by Coctables/Phenols	SCMS											
Acid Extra Prepared I		< 1.38	U	μg/l	7.04	1.38	1	SW846 8270D	24-Mar-17	25-Mar-17	MSL	1705073	
Acid Extra Prepared I 08-95-2 Surrogate re	ctables/Phenols by method SW846 3510C Phenol ecoveries:	< 1.38	U	µg/l			1	SW846 8270D	24-Mar-17	25-Mar-17	MSL	1705073	
Acid Extra Prepared I 08-95-2 Surrogate re	ctables/Phenols by method SW846 3510C Phenol ecoveries: 2-Fluorophenol	< 1.38	U	μg/l	15-11	0 %	1	SW846 8270D	24-Mar-17	25-Mar-17	"	1705073	
Acid Extra Prepared I 08-95-2 Surrogate r 167-12-4	ctables/Phenols by method SW846 3510C Phenol ecoveries: 2-Fluorophenol Phenol-d5	< 1.38	U	µg/l		0 %	1	SW846 8270D	24-Mar-17	25-Mar-17 " "	MSL "	1705073	
Acid Extra Prepared I 08-95-2 Surrogate re 167-12-4 165-62-2 SVOCs by	ctables/Phenols by method SW846 3510C Phenol ecoveries: 2-Fluorophenol Phenol-d5	< 1.38 45 31	U		15-11 15-11	0 %			24-Mar-17	" "	"	1705073	
Acid Extra Prepared I 08-95-2 Surrogate ro 167-12-4 1165-62-2 SVOCs by	ctables/Phenols by method SW846 3510C Phenol ecoveries: 2-Fluorophenol Phenol-d5 / SIM Acenaphthene	< 1.38	U	μg/l	15-11	0 %	1	SW846 8270D " " SW846 8270D SIM	24-Mar-17	25-Mar-17 " " 25-Mar-17	"	1705073	
Acid Extra Prepared I 08-95-2 Surrogate ro 667-12-4 1165-62-2 SVOCs by 3-32-9	ctables/Phenols by method SW846 3510C Phenol ecoveries: 2-Fluorophenol Phenol-d5	< 1.38 45 31	U		15-11 15-11	0 %		" " SW846 8270D	24-Mar-17	" "	"	1705073	
Acid Extra Prepared I 08-95-2 Surrogate r 67-12-4 165-62-2 SVOCs by 3-32-9	ctables/Phenols by method SW846 3510C Phenol ecoveries: 2-Fluorophenol Phenol-d5 / SIM Acenaphthene	< 1.38 45 31 < 0.070	U	μg/l	15-11 15-11 0.070	0 %		" " SW846 8270D	" "	" "	"	"""""""""""""""""""""""""""""""""""""""	
Acid Extra Prepared I 08-95-2 Surrogate ro 67-12-4 165-62-2 SVOCs by 3-32-9 08-96-8 20-12-7	ctables/Phenols by method SW846 3510C Phenol ecoveries: 2-Fluorophenol Phenol-d5 / SIM Acenaphthene Acenaphthylene	< 1.38 45 31 < 0.070 < 0.070	U	hā\J	15-11 15-11 0.070 0.070	0 % 0 % 0.043 0.044	1	" " SW846 8270D	" "	" "	"	"""""""""""""""""""""""""""""""""""""""	
Coid Extra Prepared I 08-95-2 Surrogate ro 67-12-4 165-62-2 SVOCs by 3-32-9 08-96-8 20-12-7 6-55-3	ctables/Phenols by method SW846 3510C Phenol ecoveries: 2-Fluorophenol Phenol-d5 / SIM Acenaphthene Acenaphthylene Anthracene	< 1.38 45 31 < 0.070 < 0.070 < 0.070	U	μg/l μg/l μg/l	15-11 15-11 0.070 0.070 0.070	0 % 0 % 0.043 0.044 0.037	1 1 1	" " SW846 8270D	" "	" "	"	1705073	
Coid Extra Prepared I 08-95-2 Currogate ro 67-12-4 165-62-2 CVOCs by 3-32-9 08-96-8 20-12-7 6-55-3 0-32-8 05-99-2	ctables/Phenols by method SW846 3510C Phenol ecoveries: 2-Fluorophenol Phenol-d5 / SIM Acenaphthene Acenaphthylene Anthracene Benzo (a) anthracene	< 1.38 45 31 < 0.070 < 0.070 < 0.070 < 0.070	U	hā\l hā\l	15-11 15-11 0.070 0.070 0.070 0.070	0 % 0 % 0.043 0.044 0.037 0.033	1 1 1	" " SW846 8270D	" "	" "	"	1705073	
Acid Extra Prepared I 08-95-2 Surrogate re 67-12-4 165-62-2 SVOCs by 3-32-9 08-96-8 20-12-7 6-55-3 0-32-8 05-99-2 91-24-2	ctables/Phenols by method SW846 3510C Phenol ecoveries: 2-Fluorophenol Phenol-d5 / SIM Acenaphthene Acenaphthylene Anthracene Benzo (a) anthracene Benzo (b) fluoranthene Benzo (g,h,i) perylene	< 1.38 45 31 < 0.070 < 0.070 < 0.070 < 0.070 < 0.070 < 0.070	U	hā\I hā\I hā\I	15-11 15-11 0.070 0.070 0.070 0.070 0.070	0 % 0 0 % 0.043 0.044 0.037 0.033 0.051	1 1 1 1	" " SW846 8270D	" "	" "	MSL	" " " " " " " " " " " " " " " " " " " "	
Acid Extra Prepared I 08-95-2 Surrogate ri 67-12-4 165-62-2 SVOCs by 3-32-9 08-96-8 20-12-7 6-55-3 0-32-8 05-99-2 91-24-2 07-08-9	ctables/Phenols by method SW846 3510C Phenol ecoveries: 2-Fluorophenol Phenol-d5 / SIM Acenaphthene Acenaphthylene Anthracene Benzo (a) anthracene Benzo (b) fluoranthene	< 1.38 45 31 < 0.070 < 0.070 < 0.070 < 0.070 < 0.070 < 0.070 < 0.070	U	hā\I hā\I hā\I hā\I	15-11 15-11 0.070 0.070 0.070 0.070 0.070 0.070	0 % 0 % 0.043 0.044 0.037 0.033 0.051 0.049	1 1 1 1 1	" " SW846 8270D	" "	" "	MSL	" " " " " " " " " " " " " " " " " " " "	
Acid Extra Prepared I 08-95-2 Surrogate ra 167-12-4 165-62-2 SVOCs by 3-32-9 08-96-8 20-12-7 6-55-3 0-32-8 05-99-2 91-24-2 07-08-9 18-01-9	ctables/Phenols by method SW846 3510C Phenol ecoveries: 2-Fluorophenol Phenol-d5 / SIM Acenaphthene Acenaphthylene Anthracene Benzo (a) anthracene Benzo (b) fluoranthene Benzo (g,h,i) perylene Benzo (k) fluoranthene Chrysene	< 1.38 45 31 < 0.070 < 0.070 < 0.070 < 0.070 < 0.070 < 0.070 < 0.070 < 0.070 < 0.070 < 0.070	U	hā\I hā\I hā\I hā\I hā\I	15-11 15-11 0.070 0.070 0.070 0.070 0.070 0.070 0.070 0.070	0 % 0 % 0.043 0.044 0.037 0.033 0.051 0.049 0.038 0.039	1 1 1 1 1 1 1	" " SW846 8270D	" "	" "	" " " " " " " " " " " " " " " " " " "	" " " " " " " " " " " " " " " " " " " "	
Coid Extra Prepared I 08-95-2 Currogate ro 67-12-4 165-62-2 CVOCs by 3-32-9 08-96-8 20-12-7 6-55-3 0-32-8 05-99-2 91-24-2 07-08-9 18-01-9 3-70-3	ctables/Phenols by method SW846 3510C Phenol ecoveries: 2-Fluorophenol Phenol-d5 / SIM Acenaphthene Acenaphthylene Anthracene Benzo (a) anthracene Benzo (a) pyrene Benzo (b) fluoranthene Benzo (g,h,i) perylene Benzo (k) fluoranthene	< 1.38 45 31 < 0.070 < 0.070 < 0.070 < 0.070 < 0.070 < 0.070 < 0.070 < 0.070 < 0.070	U	hā\I hā\I hā\I hā\I hā\I	15-11 15-11 0.070 0.070 0.070 0.070 0.070 0.070 0.070	0 % 0 0 % 0.044 0.037 0.033 0.051 0.049 0.038	1 1 1 1 1 1	" " SW846 8270D	" "	" "	MSL	" " " " " " " " " " " " " " " " " " " "	
Coid Extra Prepared I 08-95-2 Surrogate ro 67-12-4 165-62-2 SVOCs by 3-32-9 08-96-8 20-12-7 6-55-3 0-32-8 05-99-2 91-24-2 07-08-9 18-01-9 3-70-3 06-44-0	ctables/Phenols by method SW846 3510C Phenol ecoveries: 2-Fluorophenol Phenol-d5 / SIM Acenaphthene Acenaphthylene Anthracene Benzo (a) anthracene Benzo (a) pyrene Benzo (b) fluoranthene Benzo (g,h,i) perylene Benzo (k) fluoranthene Chrysene Dibenzo (a,h) anthracene Fluoranthene	< 1.38 45 31 < 0.070 < 0.070 < 0.070 < 0.070 < 0.070 < 0.070 < 0.070 < 0.070 < 0.070 < 0.070 < 0.070 < 0.070	U	hā\l hā\l hā\l hā\l hā\l	15-11 15-11 0.070 0.070 0.070 0.070 0.070 0.070 0.070 0.070 0.070 0.070	0 % 0.043 0.044 0.037 0.033 0.051 0.049 0.038 0.039 0.033 0.036 0.028	1 1 1 1 1 1 1 1 1	" " SW846 8270D	" "	" "	" " " " " " " " " " " " " " " " " " "	" " " " " " " " " " " " " " " " " " " "	
Coid Extra Prepared I 08-95-2 Surrogate ri 67-12-4 165-62-2 SVOCs by 3-32-9 08-96-8 20-12-7 6-55-3 0-32-8 05-99-2 91-24-2 07-08-9 18-01-9 3-70-3 06-44-0 6-73-7	ctables/Phenols by method SW846 3510C Phenol ecoveries: 2-Fluorophenol Phenol-d5 / SIM Acenaphthene Acenaphthylene Anthracene Benzo (a) anthracene Benzo (b) fluoranthene Benzo (g,h,i) perylene Benzo (k) fluoranthene Chrysene Dibenzo (a,h) anthracene Fluoranthene Fluorene	< 1.38 45 31 < 0.070 < 0.070 < 0.070 < 0.070 < 0.070 < 0.070 < 0.070 < 0.070 < 0.070 < 0.070 < 0.070 < 0.070 < 0.070 < 0.070	U	hā\l hā\l hā\l hā\l hā\l hā\l	15-11 15-11 0.070 0.070 0.070 0.070 0.070 0.070 0.070 0.070 0.070 0.070	0 % 0 % 0.043 0.044 0.037 0.033 0.051 0.049 0.038 0.039 0.033 0.036 0.028	1 1 1 1 1 1 1 1 1	" " SW846 8270D	" "	" "	" " " " " " " " " " " " " " " " " " "	" " " " " " " " " " " " " " " " " " " "	
Acid Extra Prepared I 08-95-2 Surrogate ro 67-12-4 1165-62-2 SVOCs by 3-32-9 08-96-8 20-12-7 6-55-3 0-32-8 05-99-2 91-24-2 07-08-9 18-01-9 3-70-3 06-44-0 6-73-7 93-39-5	ctables/Phenols by method SW846 3510C Phenol ecoveries: 2-Fluorophenol Phenol-d5 / SIM Acenaphthene Acenaphthylene Anthracene Benzo (a) anthracene Benzo (b) fluoranthene Benzo (g,h,i) perylene Benzo (k) fluoranthene Chrysene Dibenzo (a,h) anthracene Fluoranthene Fluorene Indeno (1,2,3-cd) pyrene	< 1.38 45 31 < 0.070 < 0.070 < 0.070 < 0.070 < 0.070 < 0.070 < 0.070 < 0.070 < 0.070 < 0.070 < 0.070 < 0.070 < 0.070 < 0.070 < 0.070 < 0.070	U	hā\l hā\l hā\l hā\l hā\l hā\l	15-11 15-11 0.070 0.070 0.070 0.070 0.070 0.070 0.070 0.070 0.070 0.070 0.070	0 % 0 0 % 0 0 0 0 0 0 0 0 0 0 0 0 0 0 0	1 1 1 1 1 1 1 1 1 1	" " SW846 8270D	" "	" "	" " " " " " " " " " " " " " " " " " "	" " " " " " " " " " " " " " " " " " " "	
Acid Extra Prepared I 08-95-2 Surrogate ro 167-12-4 1165-62-2 SVOCs by 13-32-9 108-96-8 20-12-7 16-55-3 10-32-8 105-99-2 191-24-2 107-08-9 118-01-9 13-70-3 106-44-0 16-73-7 193-39-5 111-20-3	ctables/Phenols by method SW846 3510C Phenol ecoveries: 2-Fluorophenol Phenol-d5 / SIM Acenaphthene Acenaphthylene Anthracene Benzo (a) anthracene Benzo (a) pyrene Benzo (b) fluoranthene Benzo (g,h,i) perylene Benzo (k) fluoranthene Chrysene Dibenzo (a,h) anthracene Fluoranthene Fluorene Indeno (1,2,3-cd) pyrene Naphthalene	< 1.38 45 31 < 0.070 < 0.070 < 0.070 < 0.070 < 0.070 < 0.070 < 0.070 < 0.070 < 0.070 < 0.070 < 0.070 < 0.070 < 0.070 < 0.070 < 0.070 < 0.070 < 0.070 < 0.070	U	hā\l hā\l hā\l hā\l hā\l hā\l	15-11 15-11 0.070 0.070 0.070 0.070 0.070 0.070 0.070 0.070 0.070 0.070 0.070	0 % 0 0 % 0 0 0 0 0 0 0 0 0 0 0 0 0 0 0	1 1 1 1 1 1 1 1 1 1 1	" " SW846 8270D	" "	" "	" " " " " " " " " " " " " " " " " " "	" " " " " " " " " " " " " " " " " " " "	
Acid Extra Prepared I 08-95-2 Surrogate ri 167-12-4 165-62-2 SVOCs by 3-32-9 08-96-8 20-12-7 6-55-3 0-32-8 05-99-2 91-24-2 07-08-9 18-01-9 3-70-3 06-44-0 6-73-7 93-39-5 1-20-3 5-01-8	ctables/Phenols by method SW846 3510C Phenol ecoveries: 2-Fluorophenol Phenol-d5 / SIM Acenaphthene Acenaphthylene Anthracene Benzo (a) anthracene Benzo (a) pyrene Benzo (b) fluoranthene Benzo (g,h,i) perylene Benzo (k) fluoranthene Chrysene Dibenzo (a,h) anthracene Fluoranthene Fluorene Indeno (1,2,3-cd) pyrene Naphthalene Phenanthrene	< 1.38 45 31 < 0.070 < 0.070 < 0.070 < 0.070 < 0.070 < 0.070 < 0.070 < 0.070 < 0.070 < 0.070 < 0.070 < 0.070 < 0.070 < 0.070 < 0.070 < 0.070 < 0.070 < 0.070 < 0.070	U	ha\l ha\l ha\l ha\l ha\l ha\l ha\l ha\l	15-11 15-11 0.070 0.070 0.070 0.070 0.070 0.070 0.070 0.070 0.070 0.070 0.070 0.070	0 % 0 0 % 0 0 0 0 0 0 0 0 0 0 0 0 0 0 0	1 1 1 1 1 1 1 1 1 1 1 1 1 1 1 1 1 1 1	" " " " " " " " " " " " " " " " " " "	" "	" "	" " " " " " " " " " " " " " " " " " "		
Acid Extra Prepared I 08-95-2 Surrogate ro 167-12-4 1165-62-2 SVOCs by 13-32-9 108-96-8 20-12-7 16-55-3 10-32-8 105-99-2 91-24-2 107-08-9 118-01-9 13-70-3 106-44-0 16-73-7 93-39-5 11-20-3 15-01-8	ctables/Phenols by method SW846 3510C Phenol ecoveries: 2-Fluorophenol Phenol-d5 / SIM Acenaphthene Acenaphthylene Anthracene Benzo (a) anthracene Benzo (a) pyrene Benzo (b) fluoranthene Benzo (g,h,i) perylene Benzo (k) fluoranthene Chrysene Dibenzo (a,h) anthracene Fluoranthene Fluorene Indeno (1,2,3-cd) pyrene Naphthalene	< 1.38 45 31 < 0.070 < 0.070 < 0.070 < 0.070 < 0.070 < 0.070 < 0.070 < 0.070 < 0.070 < 0.070 < 0.070 < 0.070 < 0.070 < 0.070 < 0.070 < 0.070 < 0.070 < 0.070	U	ha\l ha\l ha\l ha\l ha\l ha\l ha\l ha\l	15-11 15-11 0.070 0.070 0.070 0.070 0.070 0.070 0.070 0.070 0.070 0.070 0.070	0 % 0 0 % 0 0 0 0 0 0 0 0 0 0 0 0 0 0 0	1 1 1 1 1 1 1 1 1 1 1	" " SW846 8270D	" "	" "	" " " " " " " " " " " " " " " " " " "	" " " " " " " " " " " " " " " " " " " "	
Acid Extra Prepared I 108-95-2 Surrogate ro 167-12-4 1165-62-2 SVOCs by 13-32-9 108-96-8 120-12-7 16-55-3 160-32-8 105-99-2 191-24-2 107-08-9 118-01-9 13-70-3 106-44-0 16-73-7 193-39-5 11-20-3 15-01-8 129-00-0 Surrogate ro	ctables/Phenols by method SW846 3510C Phenol ecoveries: 2-Fluorophenol Phenol-d5 / SIM Acenaphthene Acenaphthylene Anthracene Benzo (a) anthracene Benzo (a) pyrene Benzo (b) fluoranthene Benzo (g,h,i) perylene Benzo (k) fluoranthene Chrysene Dibenzo (a,h) anthracene Fluoranthene Fluorene Indeno (1,2,3-cd) pyrene Naphthalene Phenanthrene Pyrene	< 1.38 45 31 < 0.070 < 0.070 < 0.070 < 0.070 < 0.070 < 0.070 < 0.070 < 0.070 < 0.070 < 0.070 < 0.070 < 0.070 < 0.070 < 0.070 < 0.070 < 0.070 < 0.070 < 0.070 < 0.070	U	ha\l ha\l ha\l ha\l ha\l ha\l ha\l ha\l	15-11 15-11 0.070 0.070 0.070 0.070 0.070 0.070 0.070 0.070 0.070 0.070 0.070 0.070	0 % 0 0 % 0 0 0 0 0 0 0 0 0 0 0 0 0 0 0	1 1 1 1 1 1 1 1 1 1 1 1 1 1 1 1 1 1 1	" " " " " " " " " " " " " " " " " " "	" "	" "	" " " " " " " " " " " " " " " " " " "		

Sample Id Outfall 00 SC32732-				Client P Gulf C		:	<u>Matrix</u> Surface W		ection Date -Mar-17 10			<u>ceived</u> Mar-17	
CAS No.	Analyte(s)	Result	Flag	Units	*RDL	MDL	Dilution	Method Ref.	Prepared	Analyzed	Analyst	Batch	Cert
	als by EPA 200/6000 Series by method General Prep												
·	Preservation	Field Preserved; pH<2 confirmed		N/A			1	EPA 200/6000 methods	24-Mar-17		BK	1705119	
Total Meta	als by EPA 200 Series Meth	ods											
7440-43-9	Cadmium	0.0004	J	mg/l	0.0025	0.0004	1	EPA 200.7	27-Mar-17	28-Mar-17	EDT	1705143	X
7440-47-3	Chromium	0.00979	R01, D	mg/l	0.00500	0.00306	10	EPA 200.8	"	28-Mar-17	edt	1705144	X
7440-50-8	Copper	0.0092		mg/l	0.0050	0.0029	1	EPA 200.7	"	31-Mar-17	tbc	1705143	Х
7440-02-0	Nickel	0.0044	J	mg/l	0.0050	0.0010	1	"	"	28-Mar-17	"	"	Χ
7439-92-1	Lead	0.0095		mg/l	0.0075	0.0034	1	"	"	"	ıı	"	Χ
7440-66-6	Zinc	0.0393		mg/l	0.0050	0.0027	1	"	"	"	"	"	Χ
General C	hemistry Parameters												
7782-50-5	Total Residual Chlorine	0.079		mg/l	0.020	0.006	1	SM4500-CI-G (11)	24-Mar-17 09:36	24-Mar-17 16:41	RLT	1705084	X
Prepared	by method SM4500-NH3 Ammonia as N	0.420		mg/l	0.200	0.118	1	SM4500-NH3 C. (11)	27-Mar-17	27-Mar-17	AHK	1705173	X
	рН	7.21		pH Units			1	ASTM D 1293-99B	23-Mar-17 17:56	23-Mar-17 18:30	BD	1705079	X
	Salinity	1.70		ppt (1000)	1.00	0.144	1	SM 2520 (01)		28-Mar-17	BD	1705216	j
	Total Solids	1,940		mg/l	5.00	1.53	1	SM2540 B (11)	25-Mar-17	28-Mar-17	CMB	1705113	
	Total Suspended Solids	25.0		mg/l	2.5	0.8	1	SM2540D (11)	28-Mar-17	30-Mar-17	СМВ	1705207	X
	Total Organic Carbon	5.58		mg/l	1.00	0.246	1	SM5310B (00, 11)	30-Mar-17	30-Mar-17	RLT	1705454	×
Microbiolo	ogical Analyses												
	Fecal Coliforms	2	D	CFU/100 ml			2	SM 9222D-97	23-Mar-17 15:52	23-Mar-17 15:52	NV	1705063	X
Aquatic To Prepared	oxicity by method NA												
Analysis pe	erformed by GZA Geoenviro	nmental, Inc M	lanchester,	. CT* -									
	Aquatic Toxicity	See Report		N/A			1	EPA-821-R-02-0 12		10-Apr-17		'[none]'	
	cted Analyses by method 380670												
	erformed by Phoenix Enviro	nmental Labs, In	c. * - MAC	CT007									
. 1	Oil and Grease by EPA 1664A	< 1.4		mg/L	1.4	1.4	1	E1664A		28-Mar-17 07:58	CT007	380670A	ı

Volatile Organic Compounds - Quality Control

Analyte(s)	Result	Flag	Units	*RDL	Spike Level	Source Result	%REC	%REC Limits	RPD	RPD Limit
SW846 8260C										
atch 1705104 - SW846 5030 Water MS										
Blank (1705104-BLK1)					Pre	epared & Ar	nalyzed: 24-	Mar-17		
Benzene	< 1.00		μg/l	1.00						
Benzene	< 1.0		μg/l	1.0						
Ethylbenzene	< 1.00		μg/l	1.00						
Methyl tert-butyl ether	< 1.00		μg/l	1.00						
Naphthalene	< 1.00		μg/l	1.00						
Ethylbenzene	< 1.0		μg/l	1.0						
Toluene	< 1.00		μg/l	1.00						
Naphthalene	< 1.0		μg/l	1.0						
Toluene	< 1.0		μg/l	1.0						
Vinyl chloride	< 1.00		μg/l	1.00						
m,p-Xylene	< 2.00		μg/l	2.00						
o-Xylene	< 1.00		μg/l	1.00						
Tert-Butanol / butyl alcohol	< 10.0		μg/l	10.0						
m,p-Xylene	< 2.0		μg/l	2.0						
o-Xylene	< 1.0		μg/l	1.0						
Ethanol	< 200		μg/l	200						
Surrogate: 4-Bromofluorobenzene	50.7		μg/l		50.0		101	70-130		
Surrogate: Toluene-d8	50.3		μg/l		50.0		101	70-130		
Surrogate: 1,2-Dichloroethane-d4	51.1		μg/l		50.0		102	70-130		
Surrogate: Dibromofluoromethane	49.7		μg/l		50.0		99	70-130		
Surrogate: 4-Bromofluorobenzene	50.7		μg/l		50.0		101	70-130		
Surrogate: Toluene-d8	50.3		μg/l		50.0		101	70-130		
Surrogate: 1,2-Dichloroethane-d4	51.1		μg/l		50.0		102	70-130		
Surrogate: Dibromofluoromethane	49.7		μg/l		50.0		99	70-130 70-130		
_	43.7		р9/1			anarad 9 A				
LCS (1705104-BS1)	22.2		//			epareu & Ar	nalyzed: 24-			
Benzene	22.3		μg/l		20.0		111	70-130		
Benzene	22.3		μg/l		20.0		111	70-130		
Ethylbenzene Methyl tort but diethor	23.0		μg/l		20.0		115	70-130 70-130		
Methyl tert-butyl ether	19.9		μg/l		20.0		100			
Ethylbenzene	23.0		μg/l		20.0		115	70-130		
Naphthalene	18.6		μg/l		20.0		93	70-130		
Toluene	22.2		μg/l		20.0		111	70-130		
Naphthalene	18.6		μg/l		20.0		93	70-130		
Toluene	22.2		μg/l		20.0		111	70-130 70-130		
Vinyl chloride	21.3		μg/l		20.0		106 116	70-130 70-130		
m,p-Xylene	23.2		μg/l		20.0		116	70-130 70-130		
o-Xylene	22.7		μg/l		20.0		114	70-130 70-130		
Tert-Butanol / butyl alcohol	224		μg/l		200		112	70-130		
m,p-Xylene	23.2		μg/l		20.0		116	70-130 70-130		
o-Xylene	22.7	003	μg/l		20.0		114	70-130		
Ethanol	551	QC2	μg/l		400		138	70-130		
Surrogate: 4-Bromofluorobenzene	51.3		μg/l		50.0		103	70-130		
Surrogate: Toluene-d8	50.7		μg/l		50.0		101	70-130		
Surrogate: 1,2-Dichloroethane-d4	51.7		μg/l		50.0		103	70-130		
Surrogate: Dibromofluoromethane	51.1		μg/l		50.0		102	70-130		
Surrogate: 4-Bromofluorobenzene	51.3		μg/l		50.0		103	70-130		
Surrogate: Toluene-d8	50.7		μg/l		50.0		101	70-130		
Surrogate: 1,2-Dichloroethane-d4	51.7		μg/l		50.0		103	70-130		
Surrogate: Dibromofluoromethane	51.1		μg/l		50.0		102	70-130		
LCS Dup (1705104-BSD1)					Pre	epared & Ar	nalyzed: 24-	Mar-17		

Volatile Organic Compounds - Quality Control

Analyte(s)	Result	Flag	Units	*RDL	Spike Level	Source Result	%REC	%REC Limits	RPD	RPD Limit
SW846 8260C										
Batch 1705104 - SW846 5030 Water MS										
LCS Dup (1705104-BSD1)					Pre	epared & Ar	nalyzed: 24-	Mar-17		
Benzene	20.9		μg/l		20.0		105	70-130	6	20
Benzene	20.9		μg/l		20.0		105	70-130	6	20
Ethylbenzene	21.2		μg/l		20.0		106	70-130	8	20
Methyl tert-butyl ether	20.3		μg/l		20.0		101	70-130	2	20
Ethylbenzene	21.2		μg/l		20.0		106	70-130	8	20
Naphthalene	18.9		μg/l		20.0		94	70-130	2	20
Toluene	20.6		μg/l		20.0		103	70-130	8	20
Naphthalene	18.9		μg/l		20.0		94	70-130	2	20
Toluene	20.6		μg/l		20.0		103	70-130	8	20
Vinyl chloride	19.0		μg/l		20.0		95	70-130	11	20
m,p-Xylene	21.1		μg/l		20.0		106	70-130	9	20
o-Xylene	21.0		μg/l		20.0		105	70-130	8	20
Tert-Butanol / butyl alcohol	241		μg/l		200		121	70-130	7	20
m,p-Xylene	21.1		μg/l		20.0		106	70-130	9	20
o-Xylene	21.0		μg/l		20.0		105	70-130	8	20
Ethanol	534	QC2	μg/l		400		134	70-130	3	20
Surrogate: 4-Bromofluorobenzene	51.0		μg/l		50.0		102	70-130		
Surrogate: Toluene-d8	51.0		μg/l		50.0		102	70-130		
Surrogate: 1,2-Dichloroethane-d4	51.1		μg/l		50.0		102	70-130		
Surrogate: Dibromofluoromethane	50.8		μg/l		50.0		102	70-130		
Surrogate: 4-Bromofluorobenzene	51.0		μg/l		50.0		102	70-130		
Surrogate: Toluene-d8	51.0		μg/l		50.0		102	70-130		
Surrogate: 1,2-Dichloroethane-d4	51.1		μg/l		50.0		102	70-130		
Surrogate: Dibromofluoromethane	50.8		μg/l		50.0		102	70-130		

Semivolatile Organic Compounds by GCMS - Quality Control

analyte(s)	Result	Flag	Units	*RDL	Spike Level	Source Result	%REC	%REC Limits	RPD	RPD Limi
W846 8270D										
Satch 1705073 - SW846 3510C										
Blank (1705073-BLK1)					Pre	epared: 24-l	Mar-17 An	alyzed: 25-M	ar-17	
Benzoic acid	< 1.99	U	μg/l	1.99						
4-Chloro-3-methylphenol	< 1.23	U	μg/l	1.23						
2-Chlorophenol	< 1.26	U	μg/l	1.26						
2,4-Dichlorophenol	< 1.21	U	μg/l	1.21						
2,4-Dimethylphenol	< 1.41	U	μg/l	1.41						
4,6-Dinitro-2-methylphenol	< 1.87	U	μg/l	1.87						
2,4-Dinitrophenol	< 2.15	U	μg/l	2.15						
2-Methylphenol	< 1.45	U	μg/l	1.45						
3 & 4-Methylphenol	< 1.45	U	μg/l	1.45						
2-Nitrophenol	< 1.45	U	μg/l	1.45						
4-Nitrophenol	< 2.92	U	μg/l	2.92						
Pentachlorophenol	< 1.87	U	μg/l	1.87						
Phenol	< 0.983	U	μg/l	0.983						
2,4,5-Trichlorophenol	< 1.19	U	μg/l	1.19						
2,4,6-Trichlorophenol	< 1.08	U	μg/l	1.08						
Surrogate: 2-Fluorophenol	47.9		μg/l		50.0		96	15-110		
Surrogate: Phenol-d5	46.6		μg/l		50.0		93	15-110		
LCS (1705073-BS1)					Pre	epared: 24-l	Mar-17 An	alyzed: 25-M	ar-17	
Benzoic acid	39.8		μg/l	1.99	50.0		80	30-130		
4-Chloro-3-methylphenol	37.3		μg/l	1.23	50.0		75	30-130		
2-Chlorophenol	37.5		μg/l	1.26	50.0		75	30-130		
2,4-Dichlorophenol	36.8		μg/l	1.21	50.0		74	30-130		
2,4-Dimethylphenol	35.4		μg/l	1.41	50.0		71	30-130		
4,6-Dinitro-2-methylphenol	52.1		μg/l	1.87	50.0		104	30-130		
2,4-Dinitrophenol	42.4		μg/l	2.15	50.0		85	30-130		
2-Methylphenol	36.1		μg/l	1.45	50.0		72	30-130		
3 & 4-Methylphenol	38.6		μg/l	1.45	50.0		77	30-130		
2-Nitrophenol	40.4		μg/l	1.45	50.0		81	30-130		
4-Nitrophenol	50.0		μg/l	2.92	50.0		100	30-130		
Pentachlorophenol	32.7		μg/l	1.87	50.0		65	30-130		
Phenol	35.5		μg/l	0.983	50.0		71	30-130		
2,4,5-Trichlorophenol	42.0		μg/l	1.19	50.0		84	30-130		
2,4,6-Trichlorophenol	38.8		μg/l	1.08	50.0		78	30-130		
Surrogate: 2-Fluorophenol	45.2		μg/l		50.0		90	15-110		
Surrogate: Phenol-d5	48.4		μg/l		50.0		97	15-110		
•	40.4		рул			narod: 24 I			or 17	
LCS Dup (1705073-BSD1) Benzoic acid	40.4		ua/l	1.99	50.0	parcu. 24-l	81	alyzed: 25-M 30-130	<u>ar-17</u> 2	20
4-Chloro-3-methylphenol			μg/l	1.99	50.0		81 77	30-130 30-130	3	20
2-Chlorophenol	38.3 37.2		μg/l	1.23	50.0		7 <i>1</i> 74	30-130	o.8	20
2,4-Dichlorophenol	37.2 39.2		μg/l μg/l	1.20	50.0		74 78	30-130	6	20
2,4-Dimethylphenol	39.2 35.2			1.41	50.0		76 70	30-130	0.6	20
4,6-Dinitro-2-methylphenol	35.2 51.9		μg/l μg/l	1. 4 1 1.87	50.0		70 104	30-130	0.6	20
2,4-Dinitrophenol	51.9 45.2		μg/l μg/l	2.15	50.0		90	30-130	0.3 7	20
2-Methylphenol	45.2 34.5		μg/l μg/l	2.15 1.45	50.0		90 69	30-130	4	20
3 & 4-Methylphenol	34.5 37.6			1.45 1.45	50.0		75	30-130	3	20
2-Nitrophenol			μg/l	1.45 1.45	50.0		75 77	30-130	3 5	20
4-Nitrophenol	38.5 51.1		μg/l	2.92	50.0		102	30-130	2	20
4-Nitrophenol Pentachlorophenol	33.3		μg/l μg/l	2.92 1.87	50.0		67	30-130	2	20
Phenol	33.3 34.2			0.983	50.0		68	30-130		20
2,4,5-Trichlorophenol	34.2 44.2		μg/l μg/l	0.983 1.19	50.0		68 88	30-130 30-130	4 5	20

Semivolatile Organic Compounds by GCMS - Quality Control

Analyte(s)	Result	Flag	Units	*RDL	Spike Level	Source Result	%REC	%REC Limits	RPD	RPE Limi
SW846 8270D										
Batch 1705073 - SW846 3510C										
LCS Dup (1705073-BSD1)					Pre	epared: 24-	Mar-17 An	alyzed: 25-M	lar-17	
2,4,6-Trichlorophenol	40.4		μg/l	1.08	50.0	•	81	30-130	4	20
Surrogate: 2-Fluorophenol	40.8		μg/l		50.0		82	15-110		
Surrogate: Phenol-d5	46.0		μg/l		50.0		92	15-110 15-110		
SW846 8270D SIM			P-3··							
Batch 1705073 - SW846 3510C										
					Dr	aparad: 24	Mor 17 An	alyzed: 25-M	lor 17	
Blank (1705073-BLK2) Acenaphthene	< 0.050		ua/l	0.050	<u> </u>	epareu. 24-	iviai-17 Ali	aiyzeu. 25-iv	<u>iai-17</u>	
·			μg/l							
Acenaphthylene	< 0.050		μg/l	0.050						
Anthracene	< 0.050		μg/l	0.050						
Benzo (a) anthracene	< 0.050		μg/l	0.050						
Benzo (a) pyrene	< 0.050		μg/l	0.050						
Benzo (b) fluoranthene	< 0.050		μg/l	0.050						
Benzo (g,h,i) perylene	< 0.050		μg/l	0.050						
Benzo (k) fluoranthene	< 0.050		μg/l	0.050						
Chrysene	< 0.050		μg/l	0.050						
Dibenzo (a,h) anthracene	< 0.050		μg/l	0.050						
Fluoranthene	< 0.050		μg/l	0.050						
Fluorene	< 0.050		μg/l	0.050						
Indeno (1,2,3-cd) pyrene	< 0.050		μg/l	0.050						
Naphthalene	< 0.050		μg/l	0.050						
Phenanthrene	< 0.050		μg/l	0.050						
Pyrene	< 0.050		μg/l	0.050						
Surrogate: Benzo (e) pyrene-d12	1.28		μg/l		1.00		128	30-130		
LCS (1705073-BS2)					Pre	epared: 24-l	Mar-17 An	alyzed: 25-M	lar-17	
Acenaphthene	0.785		μg/l	0.050	1.00		78	40-140		
Acenaphthylene	0.794		μg/l	0.050	1.00		79	40-140		
Anthracene	0.665		μg/l	0.050	1.00		66	40-140		
Benzo (a) anthracene	0.740		μg/l	0.050	1.00		74	40-140		
Benzo (a) pyrene	0.822		μg/l	0.050	1.00		82	40-140		
Benzo (b) fluoranthene	0.825		μg/l	0.050	1.00		82	40-140		
Benzo (g,h,i) perylene	0.763		μg/l	0.050	1.00		76	40-140		
Benzo (k) fluoranthene	0.778		μg/l	0.050	1.00		78	40-140		
Chrysene	0.736		μg/l	0.050	1.00		74	40-140		
Dibenzo (a,h) anthracene	0.879		μg/l	0.050	1.00		88	40-140		
Fluoranthene	0.655		μg/l	0.050	1.00		66	40-140		
Fluorene	0.813		μg/l	0.050	1.00		81	40-140		
Indeno (1,2,3-cd) pyrene	0.818			0.050	1.00		82	40-140		
Naphthalene	0.870		µg/l	0.050	1.00		87	40-140		
Phenanthrene			µg/l	0.050	1.00		62	40-140		
Pyrene	0.619 0.703		μg/l μg/l	0.050	1.00		70	40-140		
				0.050						
Surrogate: Benzo (e) pyrene-d12	1.14		μg/l		1.00		114	30-130		
LCS Dup (1705073-BSD2)			_			epared: 24-		alyzed: 25-N		_
Acenaphthene	0.946		μg/l	0.050	1.00		95	40-140	19	20
Acenaphthylene	0.969		μg/l	0.050	1.00		97	40-140	20	20
Anthracene	0.765		μg/l	0.050	1.00		76	40-140	14	20
Benzo (a) anthracene	0.874		μg/l	0.050	1.00		87	40-140	17	20
Benzo (a) pyrene	0.962		μg/l	0.050	1.00		96	40-140	16	20
Benzo (b) fluoranthene	0.995		μg/l	0.050	1.00		100	40-140	19	20
Benzo (g,h,i) perylene	0.921		μg/l	0.050	1.00		92	40-140	19	20

Semivolatile Organic Compounds by GCMS - Quality Control

Analyte(s)	Result	Flag	Units	*RDL	Spike Level	Source Result	%REC	%REC Limits	RPD	RPD Limit
SW846 8270D SIM										
Batch 1705073 - SW846 3510C										
LCS Dup (1705073-BSD2)					Pr	epared: 24-l	Mar-17 An	alyzed: 25-M	lar-17	
Benzo (k) fluoranthene	0.939		μg/l	0.050	1.00		94	40-140	19	20
Chrysene	0.863		μg/l	0.050	1.00		86	40-140	16	20
Dibenzo (a,h) anthracene	1.04		μg/l	0.050	1.00		104	40-140	17	20
Fluoranthene	0.787		μg/l	0.050	1.00		79	40-140	18	20
Fluorene	0.993		μg/l	0.050	1.00		99	40-140	20	20
Indeno (1,2,3-cd) pyrene	1.05	QR2	μg/l	0.050	1.00		105	40-140	25	20
Naphthalene	1.07	QR2	μg/l	0.050	1.00		107	40-140	21	20
Phenanthrene	0.756		μg/l	0.050	1.00		76	40-140	20	20
Pyrene	0.846		μg/l	0.050	1.00		85	40-140	18	20
Surrogate: Benzo (e) pyrene-d12	1.31	QM9	μg/l		1.00		131	30-130		

Total Metals by EPA 200 Series Methods - Quality Control

Analyte(s)	Result	Flag	Units	*RDL	Spike Level	Source Result	%REC	%REC Limits	RPD	RPD Limit
EPA 200.7										
Batch 1705143 - EPA 200 Series					р.		Man 47 A	l l - 00 l	4 47	
Blank (1705143-BLK1)				0.0040	Pre	epared: 27-	Mar-1/ Ar	nalyzed: 28-M	<u> 1ar-17</u>	
Nickel 	< 0.0010	U	mg/l	0.0010						
Lead	< 0.0034	U	mg/l	0.0034						
Zinc	< 0.0027	U	mg/l	0.0027						
Copper	< 0.0029	U U	mg/l	0.0029						
Cadmium	< 0.0004	U	mg/l	0.0004	_					
LCS (1705143-BS1)						epared: 27-		nalyzed: 31-M	<u> 1ar-17</u>	
Copper	1.34		mg/l	0.0029	1.25		108	85-115		
Zinc	1.32		mg/l	0.0027	1.25		105	85-115		
Lead	1.33		mg/l	0.0034	1.25		106	85-115		
Nickel	1.28		mg/l	0.0010	1.25		103	85-115		
Cadmium	1.38		mg/l	0.0004	1.25		111	85-115		
<u>Duplicate (1705143-DUP1)</u>			Source: S	C32731-01	Pre	epared: 27-	Mar-17 Ar	nalyzed: 31-M	<u> 1ar-17</u>	
Cadmium	< 0.0020	R01, U, D	mg/l	0.0020		BRL				20
Zinc	< 0.0136	R01, U, D	mg/l	0.0136		BRL				20
Nickel	< 0.0050	R01, U, D	mg/l	0.0050		BRL				20
Lead	< 0.0171	R01, U, D	mg/l	0.0171		BRL				20
Matrix Spike (1705143-MS1)			Source: S	C32731-01	Pre	epared: 27-	Mar-17 Ar	nalyzed: 31-M	1ar-17	
Lead	1.22	D	mg/l	0.0171	1.25	BRL	97.4	70-130		
Zinc	1.27	D	mg/l	0.0136	1.25	BRL	102	70-130		
Cadmium	1.25	D	mg/l	0.0020	1.25	BRL	100	70-130		
Nickel	1.20	D	mg/l	0.0050	1.25	BRL	96	70-130		
EPA 200.8										
Batch 1705144 - EPA 200 Series										
Blank (1705144-BLK1)					Pre	epared: 27-	Mar-17 Ar	nalyzed: 28-M	<u> 1ar-17</u>	
Copper	< 0.00009	U	mg/l	0.00009						
Chromium	< 0.00031	U	mg/l	0.00031						
LCS (1705144-BS1)					Pre	epared: 27-	Mar-17 Ar	nalyzed: 28-M	1ar-17	
Chromium	0.103	D	mg/l	0.00306	0.100	•	103	85-115		
Copper	0.0961	D	mg/l	0.00089	0.100		96	85-115		
Duplicate (1705144-DUP1)			•	C32731-01		enared: 27-		nalyzed: 28-M	1ar-17	
Copper	0.117	R01, R06, D	mg/l	0.00089	<u></u>	0.0972	<u> </u>	idiy2cd. 20 iv	18	20
Matrix Spike (1705144-MS1)		1,00, D	Source: S	C32731-01	<u>P</u> re	<u>epared: 2</u> 7-	<u>Mar-17</u> Ar	nalyzed: 28-M	<u>1ar-17</u>	
Copper	0.218	D	mg/l	0.00089	0.100	0.0972	121	70-130	<u></u>	
Post Spike (1705144-PS1)			-	C32731-01				nalyzed: 28-M	1ar-17	
opino (11001-FT-101)			<u> </u>				11 /1	, <u>-</u> 1V	191 11	

General Chemistry Parameters - Quality Control

Analyte(s)	Result	Flag	Units	*RDL	Spike Level	Source Result	%REC	%REC Limits	RPD	RPD Limit
ASTM D 1293-99B										
Batch 1705079 - General Preparation										
Reference (1705079-SRM1)					Pre	pared & An	alyzed: 23	8-Mar-17		
рН	6.09		pH Units		6.00		102	97.5-102. 5		
Reference (1705079-SRM2)					Pre	pared & An	alyzed: 23	s-Mar-17		
pH	5.99		pH Units		6.00		100	97.5-102.		
SM 2520 (01)								5		
Batch 1705216 - General Preparation										
Reference (1705216-SRM1)					Pre	pared & An	alvzed: 28	8-Mar-17		
Salinity	10.2		ppt (1000)	1.00	10.0	, , , , , , , , , , , , , , , , , , , 	102	90-110		
Reference (1705216-SRM2)			,		Pre	pared & An	alyzed: 28	s-Mar-17		
Salinity	10.2		ppt (1000)	1.00	10.0		102	90-110		
SM2540 B (11)										
Batch 1705113 - General Preparation										
Blank (1705113-BLK1)					Pre	pared: 25-N	Mar-17 Aı	nalyzed: 28-N	1ar-17	
Total Solids	< 5.00		mg/l	5.00						
LCS (1705113-BS1)					Pre	pared: 25-N	Mar-17 Aı	nalyzed: 28-N	<u>1ar-17</u>	
Total Solids	1080		mg/l	10.0	1100		98	90-110		
<u>Duplicate (1705113-DUP1)</u>			Source: SC:	32732-01	Pre	pared: 25-N	Mar-17 Aı	nalyzed: 28-N	<u>1ar-17</u>	
Total Solids	1930		mg/l	5.00		1940			0.3	5
SM2540D (11)										
Batch 1705207 - General Preparation										
Blank (1705207-BLK1)					Pre	pared: 28-N	Mar-17 Aı	nalyzed: 30-N	<u>1ar-17</u>	
Total Suspended Solids	< 0.5		mg/l	0.5						
LCS (1705207-BS1)					Pre	pared: 28-N	Mar-17 Ar	nalyzed: 30-N	<u>1ar-17</u>	
Total Suspended Solids	102		mg/l	10.0	100		102	90-110		
<u>SM4500-Cl-G (11)</u>										
Batch 1705084 - General Preparation										
Blank (1705084-BLK1)					<u>Pre</u>	pared & An	alyzed: 24	-Mar-17		
Total Residual Chlorine	< 0.020		mg/l	0.020						
LCS (1705084-BS1)						pared & An				
Total Residual Chlorine	0.048		mg/l	0.020	0.0500		96	90-110		
Duplicate (1705084-DUP1)			Source: SC:		Pre	pared & An	ialyzed: 24	<u>-Mar-17</u>	0.0	00
Total Residual Chlorine	0.034		mg/l	0.020	D.	0.034	-11-04	M 47	0.6	20
Matrix Spike (1705084-MS1) Total Residual Chlorine	0.056	QM9	Source: SC: mg/l	0.020	0.0500	pared & An 0.034	<u>181yzed: 24</u> 44	- <u>Mar-17</u> 80-120		
Matrix Spike Dup (1705084-MSD1)	0.030	QIVIO	Source: SC:			pared & An				
Total Residual Chlorine	0.057	QM9	mg/l	0.020	0.0500	0.034	45	80-120	0.9	200
Reference (1705084-SRM1)			9			pared & An				
Total Residual Chlorine	0.111		mg/l	0.020	0.112		99	85-115		
SM4500-NH3 C. (11)										
Batch 1705173 - General Preparation										
Blank (1705173-BLK1)					Pre	pared & An	ialyzed: 27	<u>'-Mar</u> -17		
Ammonia as N	< 0.200		mg/l	0.200	<u> </u>					
LCS (1705173-BS1)			-		<u>Pr</u> e	pared & An	ialyzed: 27	'-Mar-17		
Ammonia as N	4.90		mg/l	0.200	5.00		98	90-110		
Matrix Spike (1705173-MS1)			Source: SC:	32732-01	Pre	pared & An	alyzed: 27	'-Mar-17		

General Chemistry Parameters - Quality Control

Analyte(s)	Result	Flag	Units	*RDL	Spike Level	Source Result	%REC	%REC Limits	RPD	RPD Limit
SM4500-NH3 C. (11)										
Batch 1705173 - General Preparation										
Reference (1705173-SRM1)					Pre	epared & A	nalyzed: 27	-Mar-17		
Ammonia as N	1.82	QM9	mg/l	0.200	2.16		84	86-114		
SM5310B (00, 11)										
Batch 1705454 - General Preparation										
Blank (1705454-BLK1)					Pre	epared: 30-	Mar-17 An	nalyzed: 31-N	1ar-17	
Total Organic Carbon	< 1.00		mg/l	1.00						
LCS (1705454-BS1)					Pre	epared: 30-	Mar-17 An	nalyzed: 31-M	<u>1ar-17</u>	
Total Organic Carbon	17.0		mg/l	1.00	15.0		113	85-115		
<u>Duplicate (1705454-DUP1)</u>			Source: SO	C32731-01	Pre	epared & A	nalyzed: 30	-Mar-17		
Total Organic Carbon	2.28		mg/l	1.00		2.39			5	20
Matrix Spike (1705454-MS1)			Source: SO	C32731-01	Pre	epared & A	nalyzed: 30	-Mar-17		
Total Organic Carbon	7.98		mg/l	1.00	5.00	2.39	112	70-130		
Matrix Spike Dup (1705454-MSD1)			Source: SO	C32731-01	Pre	epared & A	nalyzed: 30	-Mar-17		
Total Organic Carbon	8.04		mg/l	1.00	5.00	2.39	113	70-130	0.8	30
Reference (1705454-SRM1)					Pre	epared: 30-	Mar-17 An	nalyzed: 31-N	1ar-17	
Total Organic Carbon	22.0		mg/l	1.00	20.0		110	85-115		

Subcontracted Analyses - Quality Control

Analyte(s)	Result	Flag	Units	*RDL	Spike Level	Source Result	%REC	%REC Limits	RPD	RPD Limit
E1664A Batch 380670A - 380670										
BLK (BX93085-BLK)			Source: SC			epared & Ar	nalyzed: 28-	<u>Mar-17</u>		
Oil and Grease by EPA 1664A LCS (BX93085-LCS)	< 1.4		mg/L Source: SO	1.4 232732-01	40 <u>Pre</u>	epared: A	nalyzed: 28	- -Mar-17		
Oil and Grease by EPA 1664A	40.20		mg/L	1.4	40		101	85-115		20

Notes and Definitions

reporting limit.

Notes an	d Definitions
D	Data reported from a dilution
J	Detected above the Method Detection Limit but below the Reporting Limit; therefore, result is an estimated concentration (CLP J-Flag).
QC2	Analyte out of acceptance range in QC spike but no reportable concentration present in sample.
QM5	The spike recovery was outside acceptance limits for the MS, MSD and/or PS due to matrix interference. The LCS and/or LCSD were within acceptance limits showing that the laboratory is in control and the data is acceptable.
QM9	The spike recovery for this QC sample is outside the established control limits. The sample results for the QC batch were accepted based on LCS/LCSD or SRM recoveries within the control limits.
QR2	The RPD result exceeded the QC control limits; however, both percent recoveries were acceptable. Sample results for the QC batch were accepted based on percent recoveries and completeness of QC data.
R01	The Reporting Limit has been raised to account for matrix interference.
R06	MRL raised to correlate to batch QC reporting limits.
U	Analyte included in the analysis, but not detected at or above the MDL.
dry	Sample results reported on a dry weight basis
NR	Not Reported
RPD	Relative Percent Difference
ClHT	The method for residual chlorine indicates that samples should be analyzed immediately. 40 CFR 136 specifies a holding time of 15 minutes from sampling to analysis. Therefore all aqueous residual chlorine samples not analyzed in the field are considered out of hold time at the time of sample receipt.
OG	The required Matrix Spike and Matrix Spike Duplicate (MS/MSD) for Oil & Grease method 1664B can only be analyzed when the client has submitted sufficient sample volume. An extra liter per MS/MSD is required to fulfill the method QC criteria. Please refer to Chain of Custody and QC Summary (MS/MSD) of the Laboratory Report to verify ample sample volume was submitted to fulfill the requirement.
рН	The method for pH does not stipulate a specific holding time other than to state that the samples should be analyzed as soon as possible. For aqueous samples the 40 CFR 136 specifies a holding time of 15 minutes from sampling to analysis. Therefore all aqueous pH samples not analyzed in the field are considered out of hold time at the time of sample receipt. All soil samples are analyzed as soon as possible after sample receipt.
LIV	The initial volume for this sample has been reduced due to sample matrix and/or historical data therefore elevating the

<u>Laboratory Control Sample (LCS)</u>: A known matrix spiked with compound(s) representative of the target analytes, which is used to document laboratory performance.

Matrix Duplicate: An intra-laboratory split sample which is used to document the precision of a method in a given sample matrix.

Matrix Spike: An aliquot of a sample spiked with a known concentration of target analyte(s). The spiking occurs prior to sample preparation and analysis. A matrix spike is used to document the bias of a method in a given sample matrix.

<u>Method Blank</u>: An analyte-free matrix to which all reagents are added in the same volumes or proportions as used in sample processing. The method blank should be carried through the complete sample preparation and analytical procedure. The method blank is used to document contamination resulting from the analytical process.

Method Detection Limit (MDL): The minimum concentration of a substance that can be measured and reported with 99% confidence that the analyte concentration is greater than zero and is determined from analysis of a sample in a given matrix type containing the analyte.

Reportable Detection Limit (RDL): The lowest concentration that can be reliably achieved within specified limits of precision and accuracy during routine laboratory operating conditions. For many analytes the RDL analyte concentration is selected as the lowest non-zero standard in the calibration curve. While the RDL is approximately 5 to 10 times the MDL, the RDL for each sample takes into account the sample volume/weight, extract/digestate volume, cleanup procedures and, if applicable, dry weight correction. Sample RDLs are highly matrix-dependent.

<u>Surrogate</u>: An organic compound which is similar to the target analyte(s) in chemical composition and behavior in the analytical process, but which is not normally found in environmental samples. These compounds are spiked into all blanks, standards, and samples prior to analysis. Percent recoveries are calculated for each surrogate.

<u>Continuing Calibration Verification:</u> The calibration relationship established during the initial calibration must be verified at periodic intervals. Concentrations, intervals, and criteria are method specific.



GEOTECHNICAL

ENVIRONMENTAL

ECOLOGICAL

WATER

CONSTRUCTION MANAGEMENT

77 Batson Drive Manchester, CT 06042 T: 860.643.9560 F: 860.646.7169



ACUTE AQUATIC TOXICITY TEST REPORT

Chelsea Sandwich, LLC Chelsea Sandwich Terminal Chelsea, MA

Test Start Date	3/24/17	
Test Period:	March 2017	

Report Prepared by:

New England Bioassay A Division of GZA GeoEnvironmental, Inc. 77 Batson Dr. Manchester, CT 06042

NEB Project Number: 05.0045458.00

Report Date: April 10, 2017

Report Submitted to:

Eurofins Spectrum Analytical, Inc. 11 Almgren Drive Agawam, MA 01001

Sample ID: <u>SC32732-01/SC32731-01</u>

If you have any questions concerning these results, please contact the Lab Manager, Kim Wills, at (860) 858-3153 or kimberly.wills@gza.com.

Whole Effluent Toxicity Testing Report Instruction Form

Client Name/Project: <u>Eurofins / Chelsea Sandwich Terminal</u> Test Date:	3/24/17
Sample ID: <u>SC32732-01/SC32731-01</u>	
Your results were as follows:	
Monitoring Only	
□ Fail – Please proceed according to the instructions in your permit.	
□ Invalid - Retesting is still required. Retest report will be sent at a later date	e under separate cover.
□ Original Test Invalid – Valid retest performed. Both test and retest results a	are attached.
☐ Retesting will be or has been performed according to the Case 1 Protocols out of EPA-New England's species-specific, self-implementing policy for alternate	
This is your case of dilution water toxicity. Please proceed a Protocols outlined in the attached copy of EPA-New England's species-specific policy for alternate dilution water. The alternate dilution water you select for should be described as follows: "synthetic laboratory water made up according protocols, by adding specified amounts of salts into deionized water in order to receiving water." Writing this letter should help you to avoid retests in the future.	pecific, self-implementing future tests for this species ing to EPA's toxicity test match the hardness of our
□ Available information is insufficient to determine whether this test passed or faile to your permit limits. Please submit a current copy of your permit to the NEB Latthe status of future tests results and help ensure your compliance with permit requ	b so that we can determine
Please complete the items on this list before reporting these results accordi	ng to the instructions in

Please complete the items on this list before reporting these results according to the instructions in the "Monitoring and Reporting" Section of your permit.

- Please complete, sign and date the upper portion of the "Whole Effluent Toxicity Test Report Certification" page which is the page directly following this page.
- Fill in the Sample Type and Sample Method (upper right) and the Permit Limits (lower left) on the New England Bioassay EPA Toxicity Test Summary Sheet(s) if they are incomplete.
- Fill in any missing information on the NEB Chain-of-Custody documents. This includes ensuring that the following information is recorded: Sampler's name and title, Facility name and address, Sample collection methods, Sample collection start and end dates and times, Types of sample, Chlorination status of samples upon shipment to NEB, Site description and Sample collection procedures.
- Monitoring results should be summarized on your monthly Discharge Monitoring Report Form.
- Signed and dated originals of this report must be submitted to the State (and Federal) Agencies specified in the "Monitoring and Reporting" section of your permit.

Questions? Please contact the Lab Manager, Kim Wills, at (860) 858-3153 or kimberly.wills@gza.com.

WHOLE EFFLUENT TOXICITY TEST REPORT CERTIFICATION (Permittee)

I certify under penalty of law that this document and all ATTACHMENTS were prepared under my direction or supervision in accordance with a system designed to assure that qualified personnel properly gather and evaluate the information submitted. Based on my inquiry of the person or persons who manage the system, or those persons who manage the system, or those persons directly responsible for gathering the information, the information submitted is, to the best of my knowledge and belief, true, accurate, and complete. I am aware that there are significant penalties for submitting false information, including the possibility of fine and imprisonment for knowing violations.

Executed on	[Date]	[Authorized Signature]
		[Print or Type Name and Title]
		[Print or Type the Permittee's Name]
		[Print or Type the NPDES Permit No.]

Since the WET test and report check is complicated, the New England Bioassay Aquatic Toxicity Laboratory has certified the validity of the WET test data in the section below. Please note that this does not relieve the permittee from its responsibility to sign and certify the report under 40 C.F.R. S 122.41(k).

WHOLE EFFLUENT TOXICITY TEST REPORT CERTIFICATION (Bioassay Laboratory)

I certify under penalty of law that this document and all ATTACHMENTS were prepared under my direction or supervision in accordance with a system designed to assure that qualified personnel properly gather and evaluate the information submitted. Based on my inquiry of the person or persons who manage the system, or those persons who manage the system, or those persons directly responsible for gathering the information, the information submitted is, to the best of my knowledge and belief, true, accurate, and complete. I am aware that there are significant penalties for submitting false information, including the possibility of fine and imprisonment for knowing violations.

Executed on

[Authorized Signature]

Kim Wills, Laboratory Manager [Print or Type Name and Title]

New England Bioassay

[Print or Type Name of Bioassay Laboratory]

24. Telephone Contacts

If you have questions, please contact Joy Hilton, Water Technical Unit, at (617) 918-1877 or David McDonald, Ecosystem Assessment Unit, at (617) 918-8609.

NEW ENGLAND BIOASSAY, A DIVISION OF GZA EPA TEST SUMMARY SHEET

Facility Name: Chelse	a Sandwich Terminal	Test Start Date:	3/24/17
NPDES Permit Numbe	r:MA0003280	Outfall Number:	001
Test Type	Test Species	Sample Type	Sample Method
XAcute	_ Fathead Minnow	_ Prechlorinated	<u>X</u> Grab
Chronic	_ Ceriodaphnia Dubia	_ Dechlorinated	_ Composite
_ Modified	_ Daphnia Pulex	Unchlorinated	_ Flow-thru
(Chronic reporting	XMysid Shrimp	Chlorinated	Other
LC50 values)	_ Sheepshead		
_ 24-Hour Screening			F-
	_Sea Urchin	TRC conc. <u>0.007</u> mg	/L
	Selenastrum		
75 11 11 17 17 1	_Other		
Dilution Water		1	.1 11 1
		ly upstream of or away from	
		Chelsea River	
		hardness to generally reflect	
Synthetic water prop	ored using either Milliners	Mill-Q or equivalent deioni	zad water and
		nbined with mineral water;	zed water and
	ixed with deionized water;		
Other			
_ 0000			
Effluent Sampling Date	e(s):3/23/17		
Effluent Concentrations	s Tested (in%):06.25	5 12.5 25 50 100	
	t Concentration):n		
(2 011110 211111			
Was effluent salinity ac	ljusted? Yes If yes, to	what value? 25 ppt	
Reference Toxicant test	t date: 3/1/17 Re	eference Toxicant Test Acce	ptable: Yes X No_
Age and Age Range of	Test Organisms 4 days (< 24 hours) Source of Organ	isms <u>NEB</u>
	TEST RESULTS &	PERMIT LIMITS	
	Test Acceptab		
		:	
A. Synthetic Water Cor	ntrol		
Mean Control Survival:	:100%N	Mean Control Reproduction:	N/A
Mean Control Weight:		Mean Control % Fertilization	
B. Receiving Water Co	ntrol		
Mean Control Survival	:100%N	Mean Control Reproduction:	N/A
Mean Control Weight:	N/AN	Mean Control % Fertilization	: <u>N/A</u>
C. Lab Culture Control	Yes_ No X		
D. Thiosulfate Control	Yes_ No X		
	Test Var	iability	
TE ADMICED (11)	NT/A		
Test PMSD (growth) Test PMSD (reproducti	N/A on.) N/A		

Permit Limits & Test Results

<u>Limits</u>		Results
LC50 N/A	LC50	>100%
	Upper Value	±∞
	Lower Value	100%
	Data Analysis	
	Method Used	Graphical
A-NOEC N/A	A-NOEC	100%
C-NOEC N/A	C-NOEC	N/A
	LOEC	N/A
IC25 N/A	IC25	
IC50 N/A	IC50	*********

PMSD Comparison Discussion - N/A

Concentration-Response Evaluation

The concentration-response relationship observed in this data set corresponds to the following item number in Chapter Four of "Method Guidance and Recommendations for Whole Effluent Toxicity (WET) Testing (40 CFR Part 136)", EPA 821-B-00-004, July 2000:

<u>X</u> 1.	Ideal concentration-response relationship
_ 2.	All or nothing response
_ 3.	Stimulatory response at low concentrations and detrimental effects at higher concentrations
_ 4.	Stimulation at low concentrations but no significant effect at higher concentrations
_ 5.	Interrupted concentration-response: significant effects bracketed by non-significant effects
_ 6.	Interrupted concentration-response: non-significant effects bracketed by significant effects
_ 7.	Significant effects only at highest concentration
_ 8.	Significant effects at all test concentrations but flat concentration-response curve
_ 9.	Significant effects at all test concentrations with a sloped concentration-response curve
_ 10.	Inverse concentration-response relationship

The concentration-response relationship was reviewed according to the above guidance document and the following determination was made:

_ 3. Results are inconclusive and the test should be repeated with a newly collected sample. An

		•
<u>X</u>	1.	Results are reliable and should be reported.
_	2.	Results are anomalous. An explanation is provided in the body of the report.

explanation is provided in the body of the report.

NEW ENGLAND BIOASSAY, A DIVISION OF GZA EPA TEST SUMMARY SHEET

Facility Name: Chelse	ea Sandwich Terminal	Test Start Date:	3/24/17
•	er: MA0003280	Outfall Number:	001
Test Type	Test Species	Sample Type	Sample Method
XAcute	_ Fathead Minnow	Prechlorinated	XGrab
Chronic	_ Ceriodaphnia Dubia	_ Dechlorinated	_Composite
Modified			- Composite
_	_ Daphnia Pulex	Unchlorinated	_ Flow-thru
(Chronic reporting		Chlorinated	_ Other
LC50 values)			
_ 24-Hour Screening			
	_Sea Urchin	TRC conc. <u>0.007</u> mg	ī/L
	Selenastrum		
	_Other		
Dilution Water			
X Receiving water col	lected at a point immediat	tely upstream of or away from	the discharge;
(Receiving water na	me and sampling location	: Chelsea River) 1
Alternate Surface W	ater of known quality and	a hardness to generally reflect	et the characteristics
of the receiving water	r: (Surface water name:		Y
Synthetic water pren	ared using either Millinor	e Mill-Q or equivalent deioni	zed water and
		ombined with mineral water;	zea water and
	ixed with deionized water		
		,	
_ Other		·	
Effluent Sampling Date	e(s): 3/23/17		
ziiidiii saiiipiiig zaii	3,23,17		
Effluent Concentration	s Tested (in%): 0 6.2	<u>25 12.5 25 50 100</u>	
	t Concentration): monit		
(1 cimit Eimi	<u>mon</u>	ornig only	
Was effluent salinity ad	djusted? Yes If yes, to	o what value? 25 ppt	
		<u>=</u> pp	
Reference Toxicant tes	t date: 3/1/17 F	Reference Toxicant Test Acce	ptable: Yes X No _
A so and A so Dones of	Took Onconients 12 down	(24 kassas) Sassas a SO	A.T
Age and Age Kange of	Test Organisms 12 days	(<24 hours) Source of Org	anisms AI
	TEST RESULTS &	&PERMIT LIMITS	
	Test Accepta	bility Criteria	
A Complement Water Con			
A. Synthetic Water Con			B T / A
Mean Control Survival		Mean Control Reproduction:	
Mean Control Weight:	N/A	Mean Control % Fertilization	: <u>N/A</u>
B. Receiving Water Co			
Mean Control Survival		Mean Control Reproduction:	
Mean Control Weight:	N/A	Mean Control % Fertilization	: <u>N/A</u>
C. Lab Culture Control	Yes_ No \underline{X}		
D. This sulfate Control	Vaa Na V		
D. Thiosulfate Control		riability.	
	1est Va	riability	
The ADMOD A STATE	3.7/4		
Test PMSD (growth)	N/A		
Test PMSD (reproducti	ion.) <u>IN/A</u>		

Permit Limits & Test Results

	Limits		Results
LC50	N/A	LC50	>100%
		Upper Value	$\pm \infty$
		Lower Value	100%
		Data Analysis	
		Method Used _	Graphical
A-NOEC_	N/A	A-NOEC	100%
C-NOEC _	N/A	C-NOEC	N/A
		LOEC	N/A
IC25	N/A	IC25	******
IC50	N/A	IC50	- Trongar

PMSD Comparison Discussion - N/A

Concentration-Response Evaluation

The concentration-response relationship observed in this data set corresponds to the following item number in Chapter Four of "Method Guidance and Recommendations for Whole Effluent Toxicity (WET) Testing (40 CFR Part 136)", EPA 821-B-00-004, July 2000:

<u>X</u> 1.	Ideal concentration-response relationship
_ 2.	All or nothing response
_ 3.	Stimulatory response at low concentrations and detrimental effects at higher concentrations
_ 4.	Stimulation at low concentrations but no significant effect at higher concentrations
	Interrupted concentration-response: significant effects bracketed by non-significant effects
_ 6.	Interrupted concentration-response: non-significant effects bracketed by significant effects
	Significant effects only at highest concentration
_ 8.	Significant effects at all test concentrations but flat concentration-response curve
_ 9.	Significant effects at all test concentrations with a sloped concentration-response curve
10.	Inverse concentration-response relationship

The concentration-response relationship was reviewed according to the above guidance document and the following determination was made:

- X 1. Results are reliable and should be reported.
- 2. Results are anomalous. An explanation is provided in the body of the report.
- 3. Results are inconclusive and the test should be repeated with a newly collected sample. An explanation is provided in the body of the report.

MYSIDOPSIS BAHIA AQUATIC TOXICITY TEST REPORT

<u>Test Reference Manual</u>: EPA 821-R-02-012, "Methods for Measuring the Acute Toxicity of

Effluents and Receiving Waters to Freshwater Organisms and

Marine Organisms", Fifth Edition

Test Method: Mysidopsis bahia Acute Toxicity Test – Method 2007.0

<u>Test Type</u>: Acute Static Non-Renewal Saltwater Test

Salinity: $25 \text{ ppt} \pm 10\%$ for all dilutions by dry ocean salts (Instant Ocean)

Temperature: 25 ± 1 °C

<u>Light Quality</u>: Ambient Laboratory Illumination

Photoperiod: 16 hours light, 8 hours dark

Test Chamber Size: 250 mL

Test Solution Volume: Minimum 200 mL

Age of Test Organisms: 4 days

Number of Mysids

Per Test Chamber: 10

Number of Replicate Test
Chambers Per Treatment: 4

Total Number of Mysids

Per Test Concentration: 40

Feeding Regime: Light feeding using concentrated *Artemia* nauplii while holding

prior to initiating the test.

Aeration: Aerated at <100 bubbles/minute

Dilution Water: Chelsea River

Alternate Control Water: NEB Artificial Salt Water (salinity 25 ppt)

Effluent Concentrations: 0%, 6.25%, 12.5%, 25%, 50% and 100% effluent

Test Duration: 48 hours

Effect measured: Mortality – no movement of body appendages on gentle prodding.

Test Acceptability: $\geq 90\%$ survival of test organisms in control solution Yes X No

<u>Sampling Requirements:</u> Samples first used within 36 hours of collection Yes \underline{X} No

Sample Volume Required: Minimum 2 liters

Test Organism Source: New England Bioassay

Test Acceptability Criteria	: Mean Alternate Water Control Mean Dilution Water Control	
Test Results:	<u>Li</u>	mits Results
	48-hour LC50 N. Upper Value Lower Value Data Analysis Method Used A-NOEC	$/A$ $ \begin{array}{r} >100\% \\ \pm \infty \\ \hline 100\% \\ \hline Graphical \\ 100\% \end{array} $
Reference Toxicant Data:	Date: Toxicant: Dilution Water: Toxicant Source: Organism Source: 48-hour LC50: In Acceptable Range:	3/1/17 Sodium Dodecyl Sulfate NEB Artificial Salt Water New England Bioassay New England Bioassay20.3 mg/L YesX No
Dechlorination Procedures	: Chlorine is measured using 4	500 CL-G DPD Colorimetric Method.
X Dechlorination was not rec	quired.	
Since dechlorination of the e	ffluent was necessary, a thiosul	the sample prior to test initiation. Ifate control of diluent water spiked Chlorine was mg/L in a
Chlorine Measurement was filtered sample.	elevated due to interference. C	Chlorine was mg/L in a
Total Residual Chlorine wa	s re-measured following aeration	on, and was found to be mg/L.
Additional Notes or Other	Conditions Affecting the Test	į.

MENIDIA BERYLLINA AQUATIC TOXICITY TEST REPORT

Test Reference Manual: EPA 821-R-02-012, "Methods for Measuring the Acute Toxicity of

Effluents and Receiving Waters to Freshwater Organisms and

Marine Organisms", Fifth Edition

<u>Test Method:</u> Menidia beryllina Acute Toxicity Test – Method 2006.0

<u>Test Type</u>: Acute Static Non-Renewal Saltwater Test

Salinity: 25 ppt \pm 2 ppt by adding dry ocean salts (Instant Ocean)

Temperature: $25 \pm 1^{\circ}$ C

Light Quality: Ambient Laboratory Illumination

Photoperiod: 16 hours light, 8 hours dark

Test Chamber Size: 250 mL

Test Solution Volume: Minimum 200 mL/replicate

Age of Test Organisms: 12 days old (24 hour age range)

Number of Fish Per

Test Chamber: 10

Number of Replicate Test
Chambers Per Treatment: 4

Total Number of Organisms
Per Test Concentration: 40

Feeding Regime: Light feeding using concentrated Artemia nauplii while holding

prior to initiating the test.

Aeration: Aerated at <100 bubbles/minute

Dilution Water: Chelsea River

Alternate Control Water: NEB Artificial Salt Water (salinity 25 ppt)

Effluent Concentrations: 0%, 6.25%, 12.5%, 25%, 50% and 100% effluent

Test Duration: 48 hours

Effect measured: Mortality – no movement on gentle prodding.

Test Acceptability: $\geq 90\%$ survival of test organisms in control solution Yes X No

<u>Sampling Requirements:</u> Samples first used within 36 hours of collection Yes \underline{X} No_

Sample Volume Required: Minimum 2 liters

Test Organism Source: Aquatic Indicators

Test Acceptability Criteria	:Mean Alternate Water Contro	
	Mean Dilution Water Control	Survival = 100%
Test Results:	<u>Li</u>	mits Results
	48-hour LC50 N Upper Value Lower Value Data Analysis Method Used A-NOEC	$7/A$ $>100\%$ $\pm \infty$ 100% $Craphical$ 100%
Reference Toxicant Data:	Date: Toxicant: Dilution Water: Toxicant Source: Organism Source: 48-hour LC50: In Acceptable Range:	3/1/17 Sodium Dodecyl Sulfate NEB Artificial Salt Water New England Bioassay Aquatic Indicators 8.49 mg/L Yes X No
Dechlorination Procedures	: Chlorine is measured using 4	500 CL-G DPD Colorimetric Method.
X Dechlorination was not re	quired.	
Since dechlorination of the e	ffluent was necessary, a thiosu	the sample prior to test initiation. Ifate control of diluent water spiked Chlorine was mg/L in a
_ Chlorine Measurement was filtered sample.	s elevated due to interference.	Chlorine was mg/L in a
_Total Residual Chlorine wa	s re-measured following aeration	on, and was found to be mg/L.
Additional Notes or Other	Conditions Affecting the Test	:
-		
<u> </u>		
F= = =================================		

NEW ENGLAND BIOASSAY ACUTE TOXICITY DATA FORM COVER SHEET FOR LC50 TESTS

CLIENT:		ctrum Analytical		M.bahia TEST ID#	17-408a
ADDRESS:		ngren Drive		M.beryllina TEST ID#	17-408b
		i, MA 01001		COC#	C37-1674/75
		Terminal Outfall 001		PROJECT #	05.0045458.00
DILUTION WATER:	Chels	sea River			
Sample Date(s):	3/	23/17	Received On:	3/24/	17
INVI	ERTEBRATES			VERTEBRATES	
TEST SE	T UP (TECH INIT)	CB		TEST SET UP (TECH INIT)	СВ
	TEST SPECIES	Mysidopsis bahia		TEST SPECIES	Menidia beryllina
	NEB LOT#	Mb17 (3-20)		NEB LOT#	Ss17AI (3-21)A
	AGE	4 days		AGE	12 days
TEST SOLUTION	ON VOLUME (mls)	200	TEST	SOLUTION VOLUME (mls)	700
NO. ORGANISMS PER	TEST CHAMBER	10	NO. ORGANI	SMS PER TEST CHAMBER	10
NO. ORGANISMS PER C	ONCENTRATION	40	NO, ORGANISI	MS PER CONCENTRATION	40
NO. ORGANISM	IS PER CONTROL	40	NO. O	RGANISMS PER CONTROL	40
	DATE	TIME		DATE	TIME
TEST START:	3/24/17	1258	TEST START:	3/24/17	1218
TEST END:	3/26/17	1300'	TEST END:	3/26/17	1248
LABORATORY CONTRO ARTIFICIAL SW:	L WATER: NEB BATCH#	CRIO37-011	Salinity (ppt)	Alkalinity (mg/L CaCO ₃₎	
RESULTS OF My	sidopsis bahia	LC50 TEST	RESULTS OF	F Menidia beryllina LC5	0 TEST
METHOD	LC50 (%)	95% Confidence Limits	METHOD	LC50 (%)	95% Confidence Limits
BINOMIAL/GRAPHICAL	>100%	100%±∞	BINOMIAL/GRAPHICAL	>100%	100%±∞
PROBIT			PROBIT		
SPEARMAN KARBER	1000/		SPEARMAN KARBER		
NOAEL	100%		NOAEL	100%	
NOEC: NO OBSERVAE	BLE EFFECT C	ONCENTRATION	I		
Comments:		243g instant oce	an was added to ~9L effluer	nt to bring salinity to 24p	pt
			ought up to 18L with D.I. w		
9			/ /	To some warming to a	y (
REVIEWD BY:			115	DATE:	4/10/17

NEW ENGLAND BIOASSAY Toxicity Test Data Sheet

NEB Test #:	17-408a	Test Organism_	Mys	idopsis bai	hia
Project #:	05.0045458.00	Organism Age:		4	days
Facility Name: _	Chelsea Sandwich Terminal	Test Duration:	48	(hours)	
Date Sampled:_	3/23/17	Beginning Date:	3/24/17	Time: _	1258
Date Received_	3/24/17	Dilution Water S	ource:	Chels	ea River
Sample ID:	Outfall 001	Salinity:	25	nı	nt

Effluent Conc. %	Number of Surviving Organisms			Dissolved Oxygen (mg/L)			Temperature (°C)				pH (su)			Salinity (ppt)	′
Initials	0	PD	PD	СВ	PD	PD	СВ	PD	PD	СВ	PD	PD	СВ	PD	PD
	0	24	48	0	24	48	0	24	48	0	24	48	0	24	48
Control A	10	10	10	7.2	5.9	5.6	24.0	25.5	24.9	8.0	7.9	7.9	25	25	25
Control B	10	10	10		5.5	5.0		25.9	25.2		7.9	7.9		25	25
Control C	10	10	10		5.6	5.1		25.9	25.1		7.9	7.9		25	25
Control D	10	10	10		5.5	5.0		25.9	25.3		7.9	7.9		25	25
Diluent A	10	10	10	7.5	5.8	4.9	24.6	25.7	25.1	7.9	7.8	7.7	25	25	25
Diluent B	10	10	10		5.8	4.5		25.7	25.2		7.8	7.6		25	25
Diluent C	10	10	10		5.6	4.3		26.0	25.3		7.7	7.6		25	25
Diluent D	10	10	10		5.4	4.4		26.0	25.3		7.7	7.6		25	25
6.25 A	10	10	10	7.4	5.6	4.8	24.6	26.0	25.2	7.8	7.8	7.7	25	25	25
6.25 B	10	10	10		5.4	4.6		25.9	25.3		7.7	7.7		25	25
6.25 C	10	10	10		5.4	4.3		25.8	25.3		7.7	7.6		25	25
6.25 D	10	10	10		5.3	4.2		25.9	25.3		7.7	7.6		25	25
12.5 A	10	10	10	7.3	5.1	4.5	24.7	25.9	25.2	7.8	7.7	7.7	25	25	25
12.5 B	10	10	10		5.2	5.1		25.9	25.1		7.8	7.8		25	25
12.5 C	10	10	10		5.7	4.9		25.6	25.2		7.8	7.7		25	25
12.5 D	10	10	10		5.6	4.8		25.7	25.3		7.8	7.7		25	25
25 A	10	10	10	7.3	5.8	5.4	24.7	25.4	25.0	7.8	7.8	7.8	25	25	25
25 B	10	10	10		5.6	4.9		25.7	25.2		7.8	7.8		25	25
25 C	10	10	10		5.5	4.7		25.7	25.3		7.8	7.8		25	25
25 D	10	10	10		5.3	4.6		25.7	25.3		7.8	7.8		25	25

LC50	Confidence Interval	A-NOEC	Computational Method		
>100%	100%±∞	100%	Graphical		

NEW ENGLAND BIOASSAY Toxicity Test Data Sheet

NEB Test #:	17-408a	Test Organism_	Mys	idopsis bah	ia
Project #:	05.0045458.00	Organism Age_		4	days
Facility Name: _	Chelsea Sandwich Terminal	Test Duration:_	48	_(hours)	
Date Sampled:	3/23/17	Beginning Date:_	3/24/17	_Time:	1258
Date Received_	3/24/17	Dilution Water S	Source	Chelsea R	River
Sample ID:	Outfall 001	Salinity.	25	pp:	t

Effluent Conc. %	Conc. Surviving			Dissolved Oxygen (mg/L)		Temperature (°C)		pH (su)			Salinity (ppt)				
Initials	0	PD	PD	СВ	PD	PD	СВ	PD	PD	СВ	PD	PD	СВ	PD	PD
	0	24	48	0	24	48	0	24	48	0	24	48	0	24	48
50 A	10	10	10	7.1	5.4	4.9	24.7	25.6	25.2	7.8	7.9	7.9	24	25	25
50 B	10	10	10		5.4	4.8		25.8	25.2		7.9	7.9		24	24
50 C	10	10	10		5.6	4.7		25.8	25.2		7.9	7.9		24	24
50 D	10	10	10		5.5	4.8		25.9	25.2		7.9	7.9		24	24
100 A	10	10	10	6.9	5.3	5.3	24.6	25.7	25.1	7.7	7.8	8.1	24	24	24
100 B	10	10	10		5.3	4.9		25.7	25.3		7.8	8.0		24	24
100 C	10	10	10		5.2	4.9		25.6	25.3		7.8	8.0		24	24
100 D	10	10	10		5.4	5.1		25.5	25.1		7.8	8.1		24	24
												,			

LC50	Confidence Interval	A-NOEC	Computational Method
>100%	100%±∞	100%	Graphical

MINIMUM REQUIRED TRIM IS TOO LARGE: 100.0, SO SK IS NOT CALCULABLE. SPEARMAN-KARBER

TRIM:

LC50: .000

95% CONFIDENCE LIMITS ARE UNRELIABLE.

.9095D-10

.00%

CONC. NUMBER NUMBER PERCENT BINOMIAL PROB.(%) EXPOSED DEAD DEAD 6.25 40. 0. .00 12.50 40. 0. .00 25.00 40. 0. .00 50.00 40. 0. .00 100.00 40. 0. .00 .9095D-10 .9095D-10 .9095D-10 .9095D-10

THE BINOMIAL TEST SHOWS THAT 100.00 AND +INFINITY CAN BE USED AS STATISTICALLY SOUND CONSERVATIVE 95 PERCENT CONFIDENCE LIMITS SINCE THE ACTUAL CONFIDENCE LEVEL ASSOCIATED WITH THESE LIMITS IS 100.0000 PERCENT. THE LC50 FOR THIS DATA SET IS GREATER THAN 100.00

WHEN THERE ARE LESS THAN TWO CONCENTRATIONS AT WHICH THE PERCENT DEAD IS BETWEEN 0 AND 100, NEITHER THE MOVING AVERAGE NOR THE PROBIT METHOD CAN GIVE ANY STATISCALLY SOUND RESULTS.

TEST NUMBER: 17-408a DURATION: 48 h DATE: 3/24/17

SAMPLE: Chelsea Sandwich SPECIES: Mysidopsis bahia

METHOD LC50 CONFIDENCE LIMITS LOWER UPPER SPAN ***** 100.000 ***** BINOMIAL MAA ***** PROBIT .000 ****** ****** ***** SPEARMAN

**** = LIMIT DOES NOT EXIST

NEW ENGLAND BIOASSAY Toxicity Test Data Sheet

NEB Test #:	17-408b	Test Organisn_	Men	idia beryllir	па
Project #:	05.0045458.00	Organism Age:	/	12	days
Facility Name:	Chelsea Sandwich Terminal	Test Duration:_	48	(hours)	
Date Sampled:	3/23/17	Beginning Date_	3/24/17	_Time:	1218
Date Received	3/24/17	Dilution Water S	Source:	Chelse	ea River
Sample ID:	Outfall 001	Salinity.	25	nn	t

Effluent Conc. %			g	Dissolved Oxygen (mg/L)		Temperature (°C)		pH (su)			Salinity (ppt)				
Initials	0	PD	PD	СВ	PD	PD	СВ	PD	PD	СВ	PD	PD	СВ	PD	PD
73-A-75	0	24	48	0	24	48	0	24	48	0	24	48	0	24	48
Control A	10	10	10	7.2	6.5	6.5	24.0	25.5	25.2	8.0	8.1	8.1	25	25	25
Control B	10	10	10		5.9	6.0		26.0	25.8		8.1	8.1		25	25
Control C	10	10	10		5.9	5.9		26.0	25.9		8.1	8.1		25	25
Control D	10	9	9		6.0	5.9		25.8	26.0		8.1	8.1		25	25
Diluent A	10	10	10	7.5	6.0	6.2	24.6	25.8	25.2	7.9	7.9	7.9	25	25	25
Diluent B	10	10	10		6.0	6.1		25.8	25.6		7.9	7.9		25	25
Diluent C	10	10	10		6.0	6.1		25.9	25.7		7.9	7.9		25	25
Diluent D	10	10	10		6.2	6.1		25.4	25.7		7.9	7.9		25	25
6.25 A	10	10	10	7.4	6.2	6.1	24.6	25.3	25.4	7.8	7.9	7.9	25	25	25
6.25 B	10	10	10		5.9	5.9		25.6	25.7		7.9	7.9		25	25
6.25 C	10	10	10		5.8	5.8		25.7	25.8		7.9	7.9		25	25
6.25 D	10	10	10		6.1	6.1		25.3	25.8		8.0	8.0		25	25
12.5 A	10	10	10	7.3	6.1	6.3	24.7	25.2	25.2	7.8	8.0	7.9	25	25	25
12.5 B	10	10	10		5.9	5.6		25.6	25.8		7.9	7.9		25	25
12.5 C	10	10	10		5.7	5.7		25.8	25.9		7.9	7.9		25	25
12.5 D	10	10	10		5.7	5.7		25.8	25.9		7.9	7.9		25	25
25 A	10	10	10	7.3	6.2	5.9	24.7	25.2	25.5	7.8	8.0	8.0	25	25	25
25 B	10	10	10		6.1	6.1		25.4	25.7		8.0	8.0		25	25
25 C	10	10	10		5.9	5.9		25.6	25.8		8.0	8.0		25	25
25 D	10	10	9		5.9	5.9		25.4	25.7		8.0	8.0		25	25

LC50	Confidence Interval	A-NOEC	Computational Method
>100%	100%±∞	100%	Graphical

NEW ENGLAND BIOASSAY Toxicity Test Data Sheet

NEB Test #: _	17-408b	Test Organisn_	Mei	nidia berylli	na
Project #:	05.0045458.00	Organism Age_		12	days
Facility Name: _	Chelsea Sandwich Terminal	Test Duration:_	48	_(hours)	
Date Sampled:_	3/23/17	Beginning Date_	3/24/17	_Time: _	1218
Date Received_	3/24/17	Dilution Water S	Source	Chelsea l	River
Sample ID	Outfall 001	Salinity	25	nr	of.

Effluent Conc. %	Conc. Surviving			Dissolved Oxygen (mg/L)		Temperature (°C)		pH (su)			Salinity (ppt)				
Initials	0	PD	PD	СВ	PD	PD	СВ	PD	PD	СВ	PD	PD	СВ	PD	PD
A NOTE TO	0	24	48	0	24	48	0	24	48	0	24	48	0	24	48
50 A	10	10	10	7.1	6.0	6.4	24.7	25.3	25.5	7.8	8.0	8.1	24	25	25
50 B	10	10	10		6.0	5.9		25.4	25.7		8.0	8.1		25	25
50 C	10	10	10		5.9	6.0		25.6	25.6		8.0	8.1		25	25
50 D	10	10	10		6.0	6.1		25.5	25.6		8.0	8.1		25	25
100 A	10	10	10	6.9	6.1	5.8	24.6	25.6	25.9	7.7	8.0	8.1	24	24	25
100 B	10	10	10		5.8	6.0		25.6	25.6		8.0	8.2		24	25
100 C	10	10	10		5.6	5.9		25.8	25.5		8.0	8.1		24	25
100 D	10	10	10		5.5	6.1		25.6	25.2		8.0	8.1		24	25

LC50	Confidence Interval	A-NOEC	Computational Method
>100%	100%±∞	100%	Graphical

CT-TOX: BINOMIAL, MOVING AVERAGE, PROBIT, AND SPEARMAN METHODS

MINIMUM REQUIRED TRIM IS TOO LARGE: 99.2, SO SK IS NOT CALCULABLE.

SPEARMAN-KARBER

TRIM:

.00%

LC50:

.000

95% CONFIDENCE LIMITS ARE UNRELIABLE.

CONC.	NUMBER	NUMBER	PERCENT	BINOMIAL	
%	EXPOSED	DEAD	DEAD	PROB.(%)	
6.25	40.	0.	.00	.9095D-10	
12.50	40.	0.	.00	.9095D-10	
25.00	40.	1.	2.50	.3729D-08	
50.00	40.	0.	.00	.9095D-10	
100.00	40.	0.	.00	.9095D-10	

THE BINOMIAL TEST SHOWS THAT 100.00 AND +INFINITY CAN BE USED AS STATISTICALLY SOUND CONSERVATIVE 95 PERCENT CONFIDENCE LIMITS SINCE THE ACTUAL CONFIDENCE LEVEL ASSOCIATED WITH THESE LIMITS IS 100.0000 PERCENT. THE LC50 FOR THIS DATA SET IS GREATER THAN 100.00

WHEN THERE ARE LESS THAN TWO CONCENTRATIONS AT WHICH THE PERCENT DEAD IS BETWEEN 0 AND 100, NEITHER THE MOVING AVERAGE NOR THE PROBIT METHOD CAN GIVE ANY STATISCALLY SOUND RESULTS.

DATE: 3/24/17 TEST NUMBER: 17-408b DURATION: 48 h

SAMPLE: Chelsea Sandwich SPECIES: Menidia beryllina

METHOD	LC50	COL	NFIDENCE LI	IMITS
		LOWER	UPPER	SPAN
BINOMIAL	*****	100.000	*****	*****
MAA	*****	*****	*****	*****
PROBIT	*****	*****	*****	*****
SPEARMAN	.000	*****	*****	*****

NOTE: MORTALITY PROPORTIONS WERE NOT MONOTONICALLY INCREASING.
ADJUSTMENTS WERE MADE PRIOR TO SPEARMAN-KARBER ESTIMATION.

**** = LIMIT DOES NOT EXIST

INITIAL CHEMISTRY INFORMATION

CLIENT: PROJECT #

Chelsea Sandwich Terminal

05.0045458.00

RECIEPT DATE	3/2	4/17
SAMPLE	Effluent	Receiving Water
COC#	C37-1674	C37-1675
Temperature (°C)	5.0	4.8
Dissolved Oxygen (mg/L)	7.9	11.0
pH (standard units)	6.8	7.8
Conductivity (µmhos/cm)	4,084	47,950
Salinity (ppt)	2	31
Hardness (as mg/L CaCO3)	430	5700
Alkalinity (as mg/L CaCO3)	80	105
TRC - DPD (mg/L)	0.007	<0.001
INITIALS	CW	CW

Additional notes:	



SUBCONTRACT ORDER

SC32732

Spectrum Analytical

SENDING LABORATORY:

Eurofins Spectrum Analytical, Inc.

II Almgren Drive

Agawam, MA 01001 Phone: (413) 789-9018

Fax: (413) 789-4076

Project Manager: Dulce Litchfield

Project: Gulf Terminal - Chelsea, MA

RECEIVING LABORATORY:

GZA Geoenvironmental, Inc. - Manchester, CT*

77 Batson Drive

Manchester, CT 06042 Phone: (860) 286-8900 Fax: (860) 242-8389

Project #:

Gulf Chelsea

PO Number:

SC32732

BILL TO:

Eurofins Spectrum Analytical, Inc.

2425 New Holland Pike

Lancaster, PA 17601

Attention: Accounts Payable accountspayable@eurofinsus.com

PO Number:

SC32732

Laboratory ID	Sample ID	Sampled	Matrix	Analysis	Due	Comments
	SC32732-01	23-Mar-17 10:00	Surface Water	Aquatic Tox	07-Apr-17 16:00	Client ID is Outfall 003/LC50
Containers Supplied:						(37-1674
Other (L)	8					(37-167

Please send notice within 24 hours of obtaining valid data, of the results of all drinking water samples that exceed any EPA or Department-established maximum contaminant level, maximum residual disinfectant level or reportable concentration. Notice should be emailed to SpectrumLabResults@EurofinsUS.com.

Please notify <u>SpectrumLabResults@EurofinsUS.com</u> immediately and prior to conducting analysis if certification is not held for the analyses requested.

Please e-mail results in electronic format to SpectrumLabResults@EurofinsUS.com.

Received Unpreserved

Released By Date Received By 3/24/11@0945

Temp

Released By

Date

Received By

Date



SUBCONTRACT ORDER

Spectrum Analytical

SC32731

SENDING LABORATORY:

Eurofins Spectrum Analytical, Inc.

11 Almgren Drive

Agawam, MA 01001

Phone: (413) 789-9018 Fax: (413) 789-4076

Project Manager: Dulce Litchfield

Project: Gulf Terminal - Chelsea, MA

RECEIVING LABORATORY:

GZA Geoenvironmental, Inc. - Manchester, CT*

77 Batson Drive

Manchester, CT 06042 Phone: (860) 286-8900

Fax: (860) 242-8389

BILL TO:

Eurofins Spectrum Analytical, Inc.

2425 New Holland Pike

Lancaster, PA 17601

Attention: Accounts Payable

accountspayable@eurofinsus.com

PO Number:

SC32731

Project #:

Gulf Chelsea

PO Number:

SC32731

Laboratory ID	Sample ID	Sampled	Matrix	Analysis	Due	Comments
- 0	SC32731-01	23-Mar-17 10:00	Surface Water	Aquatic Tox	07-Apr-17 16:00	Client ID is Chelsea Creck/LC50
ntainers Supplied						037-1675

Please send notice within 24 hours of obtaining valid data, of the results of all drinking water samples that exceed any EPA or Department-established maximum contaminant level, maximum residual disinfectant level or reportable concentration. Notice should be emailed to <u>SpectrumLabResults@EurofinsUS.com</u>.

Please notify <u>SpectrumLabResults@EurofinsUS.com</u> immediately and prior to conducting analysis if certification is not held for the analyses requested.

Please e-mail results in electronic format to SpectrumLabResults@EurofinsUS.com.

Received Unpreserved

Released By

Dute

Received By

Date

emp °(

Released By

Date

Received By

Date

NEB SALTWATER SPEC S ACCLIMATION RECORD

Species:	Client:	Quantity:	*Mortality upon arrival
Mendace berylling	Test ID:	280	, ,
	LOT#: 5517AI(3.21)A	Age:	(20)
Agraph Indicators		\$ days on 3.21-17	* Mortality > 10% - Notify management

Allowable Mortality: > 5% mortality = Notify management.

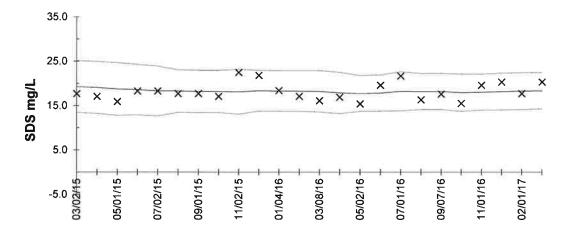
Fish = No more than 50% tank volume water change over a 12 (twelve) hour period. Allowable Acclimation:

Mysids = Need to be ± -2 ppt of test dilution water.

	Water	Water Chemistry	stry					Obser	Observations		
Date	D.O. (mg/L)	p.H. (SU)	Temp. (C) *	Alkal. (mg/L) ml titrant	Sal. (ppt) **	F	Feedings	Behavioral observations	Do organisms look stressed?	Mortalities	Comments / Treatment type
						AM	NOON PM	A = Normal, B = Erratic mov. C = Dead	Yes / No	# of dead organisms removed from tank	
3.2(-17	7,7	カン	22.5	() E	25	##	MY MB	4	2		Acclemated to ASW.
3-22-17		(23.4	[9	2	55 SP	1	2	0	Unter 1 101- Asw
3-23-17	h 2	1	22.3	1	25	à	Sy SJP	K	200	0	water D lot Asw
3-24-17	7,2		8:10	ļ	36	à	FG 538	4	2	0	Veter D lot ALL
3-26-17	6.9	{	2),6	(ر رد	200	SIP SIP	Z,	2	0	HO D 101 ASW
										200	
	1										

New England Bioassay Reference Toxicant Data: *Mysidopsis bahia* 48-hour LC50

Reference Toxicant: Sodium Dodecyl Sulfate Test Dates: March 2015 - March 2017

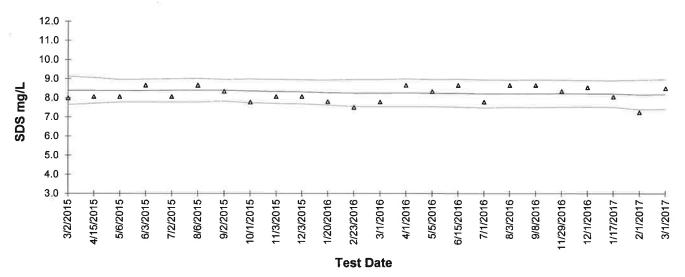


		Test Date	
×	LC50	—— Mean LC50	±2 SD

								CV National
Test ID	Date	LC ₅₀	Mean LC ₅₀	STD	-2STD	+2STD	CV	75th & 90th%
15-258	3/2/2015	17.7	19.3	2.9	13.5	25.1	0.15	0.26
15-414	4/1/2015	17.1	19.1	2.9	13.2	24.9	0.15	0.26
15-549	5/1/2015	15.9	18.7	3.0	12.8	24.7	0.16	0.26
15-704	6/1/2015	18.3	18.6	2.9	12.9	24.3	0.15	0.26
15-900	7/2/2015	18.3	18.3	2.8	12.7	23.9	0.15	0.26
15-1082	8/3/2015	17.7	18.3	2.4	13.5	23.1	0.13	0.26
15-1296	9/1/2015	17.7	18.2	2.4	13.4	23.0	0.13	0.26
15-1458	10/1/2015	17.1	18.2	2.4	13.5	23.0	0.13	0.26
15-1687	11/2/2015	22.5	18.1	2.5	13.1	23.2	0.14	0.26
15-1776	12/1/2015	21.8	18.4	2.3	13.8	23.0	0.13	0.26
16-34	1/4/2016	18.4	18.3	2.3	13.7	22.9	0.12	0.26
16-142	2/1/2016	17.1	18.3	2.3	13.7	22.8	0.12	0.26
16-338	3/8/2016	16.1	18.2	2.3	13.6	22.9	0.13	0.26
16-460	4/1/2016	16.9	17.9	2.3	13.2	22.5	0.13	0.26
16-600	5/2/2016	15.4	17.8	2.0	13.7	21.8	0.11	0.26
16-709	6/1/2016	19.6	17.9	2.0	13.8	22.0	0.11	0.26
16-849	7/1/2016	21.7	18.3	2.2	13.8	22.7	0.12	0.26
16-1058	8/1/2016	16.3	18.2	2.0	14.1	22.2	0.11	0.26
16-1256	9/7/2016	17.6	18.2	2.0	14.1	22.3	0.11	0.26
16-1471	10/5/2016	15.5	17.9	2.1	13.7	22.1	0.12	0.26
16-1590	11/1/2016	19.6	18.0	2.0	14.0	22.1	0.11	0.26
17-9	1/3/2017	20.3	18.2	2.1	14.0	22.4	0.11	0.26
17-154	2/1/2017	17.7	18.3	2.1	14.1	22.4	0.11	0.26
17-273	3/1/2017	20.3	18.4	2,1	14.3	22.5	0.11	0.26

New England Bioassay Reference Toxicant Data: *Menidia beryllina* 48-hour LC50

Reference Toxicant: Sodium Dodecyl Sulfate Test Dates: March 2015 - March 2017



r -			
Δ	LC50	Mean LC50	+/- 2 STD

								CV National	CV National
Test ID	Date	LC ₅₀	Mean LC ₅₀	STD	-2STD	+2STD	CV	75th%	90th%
15-143	3/2/2015	8.0	8.4	0.4	7.6	9.1	0.04	0.21	0.44
15-585	4/15/2015	8.1	8.4	0.3	7.7	9.1	0.04	0.21	0.44
15-623	5/6/2015	8.1	8.4	0.3	7.8	9.0	0.04	0.21	0.44
15-705	6/3/2015	8.7	8.4	0.3	7.8	9.0	0.04	0.21	0.44
15-901	7/2/2015	8.1	8.4	0.3	7.8	9.0	0.04	0.21	0.44
15-1083	8/6/2015	8.7	8.4	0.3	7.8	9.0	0.04	0.21	0.44
15-1297	9/2/2015	8.4	8.4	0.3	7.8	9.0	0.03	0.21	0.44
15-1539	10/1/2015	7.8	8.4	0.3	7.7	9.0	0.04	0.21	0.44
15-1688	11/3/2015	8.1	8.3	0.3	7.7	9.0	0.04	0.21	0.44
15-1825	12/3/2015	8.1	8.3	0.3	7.7	8.9	0.04	0.21	0.44
16-108	1/20/2016	7.8	8.3	0.3	7.6	8.9	0.04	0.21	0.44
16-260	2/23/2016	7.5	8.3	0.4	7.6	9.0	0.04	0.21	0.44
16-303	3/1/2016	7.8	8.3	0.4	7.5	9.0	0.04	0.21	0.44
16-461	4/1/2016	8.7	8.3	0.4	7.5	9.0	0.04	0.21	0.44
16-602	5/5/2016	8.3	8.3	0.4	7.5	9.0	0.04	0.21	0.44
16-798	6/15/2016	8.7	8.2	0.4	7.5	9.0	0.04	0.21	0.44
16-850	7/1/2016	7.8	8.2	0.4	7.5	8.9	0.04	0.21	0.44
16-1060	8/3/2016	8.7	8.2	0.4	7.5	8.9	0.04	0.21	0.44
16-1282	9/8/2016	8.7	8.2	0.4	7,5	8.9	0.04	0.21	0.44
16-1705	11/29/2016	8.4	8.2	0.4	7.5	8.9	0.04	0.21	0.44
16-1739	12/1/2016	8.6	8.2	0.3	7.5	8.9	0.04	0.21	0.44
17-83	1/17/2017	8.1	8.2	0.3	7.5	8.9	0.04	0.21	0.44
17-155	2/1/2017	7.3	8.2	0.4	7.4	8.9	0.05	0.21	0.44
17-278	3/1/2017	8.5	8.2	0.4	7.4	9.0	0.05	0.21	0.44

X32731 IN Special Handling:

☑ Standard TAT - 7 to 10 business days

☐ Rush TAT - Date Needed:

HANIBAL TECHNOLOGY	Featuring	SEECTRUM ANALYTICAL, INC.			01200 B0000000 21210
			 		 _

CHAIN OF CUSTODY RECORD

Project Mgr: Telephone #: Report To: Andrew Adams Chelsea, MA 02150 281 Eastern Ave Gulf Oil LP Andrew Adams 617.884.5980 Invoice To: Christopher Gill P.O No.: Wellesley, MA 02481-3705 80 William St, Suite 400 Gulf Oil LP Page ____1 of _____1 Quote/RQN: Site Name: Sampler(s): Location: Project No: All TATs subject to laboratory approval
Min. 24-hr notification needed for rushes
Samples disposed after 60 days unless otherwise instructed. 281 Eastern Ave, Chelsea Gulf Chelsea Terminal Gulf Chelsea State

MA

DI VOA Erroren Soil far Erozen	Reliperated		Ambient Ised	unbient	 	プリ	- IR	(_									
□ Present □ Intact □ Hroken	Custody Seals:		Condition upon receipt	tion upor	Fondar onder	4	9	1						7				
				-3)	Orenction Factor	8	5		375	(V)			1 - K	2	Pil	Style	
aadams@gulfoil.com, cgill@gulfoil.com	foil.com, cg	ams@gul		E-mail to:		をか	0	11:0	17	531	CU		Red	the.		Chill	the	
9			mat:	EDD format:	\$d	Temp °C		Time:		Date:			l by:	Received by:		red by:	Relinquished by:	
					3	200												
Zn - 5 μg/L												-						
Cu - 0.5 µg/L											-	-	a				•	
Cd, Pb, Ni - 0.2 ug/L		×									SW	6	1000	3-73	*	Chelsea Creek		K
Group 2 PAHs - 5 µg/L		^	×								WS	G	1000	3-73		Chelsea Creek		
Group 1 PAHs - 0.1 μg/L			×							2	WS	G	000	3-23		Chelsea Creek		
naphthalene - 5 µg/L				×							SW	G	800	3-23		Chelsea Creek		
BTEX - 2 μg/L					×					3	SW	G	1800	3.23		Chelsea Creek		
Required Minimum Levels:						×		_	-		SW	G	1000	3-73		Chelsea Creek		
* 20						×		_			SW	6	00:00	ري ري		Chelsea Creek	3/4	232
Other State-specific reporti	Char			PAHs		Amm TRC, TSS				# of '	Ma	Ту	Time:	Date:		Sample ID:	D.	Lab ID:
NJ Reduced* NJ Full*	k if a	i, Zn)			(& na			Clear (VOA		pe	6	C=Compsite		ab	_G= Grab	1/1
DQA* ASP B*	nlani-		. (Cd,		iphtha	ty, pH			r Glass	Vials	I			X3=	X2=		X1=	>
Standard No QC	o to d		Cu,		lene	, TS,			5				il Gas	Air SG=Soil Gas	A=Indoor/Ambient Air	SL=Sludge A=Iı	SO=Soil	0 =0il
čeport? □			ysis	Analysis				Containers	Cont			'ater	WW=Waste Water		SW=Surface Water	GW=Groundwater	DW=Dinking Water	DW=D
additional charges may appply	-	4	10 2	1	2	3 11		ì										
QA/QC Reporting Notes:		elow:	List Preservative Code below:	servativ	List Pre						12=		none	11:	эг 10 =Н ₃ РО ₄	7=CH3OH 8=NaHSO ₄ 9=Deionized Water 10=H ₃ PO ₄	OH 8=NaHSO	7=CH3
										cıd	6=Ascorbic Acid	6=Asc	5=NaOH	4=HNO ₃	$\mathbf{J} = \mathbf{H}_2 \mathbf{SO}_4$	$1=Na_2S2O_3$ $2=HC1$	r-rield riffered 1-1	F=Fleic

SPECTRUM ANALYTICAL, INC. HANIBAL TECHNOLOGY

CHAIN OF CUSTODY RECORD

□ Rush TAT - Date Needed:

Page __1__ of ___2

		2(5/7)	
V Standard TAT	. Spe	32 / 50	1
T - 7 to 10 business days	ecial Handling:		

							- ,						2				The second										
		J.	1/1	Relin	1			\				1	32732 0	Lab ID:		X1=	0=Oil S0=Soil	DW =Dinking Water		F=Field Filtered 1=N ₄ 7=CH3OH 8=NaHSO ₄	Project Mgr.	Telephone #:	Chelsea	281 Eastern Ave	Gulf Oil LP	Report To: Andrew Adams	SPECTR HANI
		1 Par		Religiuished by:			Outfall 003	Outfall 003	Outfall 003	Outfall 003	Outfall 003	Outfall 003	Outfall 003	Sample ID:	G= Grab		SL=Sludge A=Indoor/Ambient Air	GW=Groundwater		1=Na ₂ S2O ₃	Andrew Adams	617.884.5980	Chelsea, MA 02150	tern Ave	_p	Adams	SPECTRUM ANALYTICAL, INC. Featuring HANIBAL TECHNOLOGY
		9	Yan I	Received by:			2.23	3-73	3-73	3-23	3-73	3-23	3-23	Date:	C=Compsite	X3=	ient Air SG=Soil Gas	SW=Surface Water WW		4=HNO ₃							
		R		by:			8	060	1000	1000	8	(600	000	Time:			Gas	ww=Waste Water		5=NaOH 6=As	P.O No.:		W	80	ବ୍ର	Invoice To: Christopher Gill	
		C .	0				G S	G	G S	G S	G S	G S	G	-	ype					6=Ascorbic Acid			Wellesley, MA 02481-3705	80 William St, Suite 400	Gulf Oil LP	ıristophe	Page
	_	7	3/2	D.			WS	WS	WS	SW :	WS	WS	WS	-	voa	Viale	<i>(</i>			Acid			MA 02	1 St, Su	U	er Gill	<u> </u>
	_	7	3/	Date:			2			ω					Ambe						 -	*	481-370	ite 400			of
		9	7											-	Clear		-	Containers			Quote/RQN)5				1
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Ambient	mion up	3	E-mail to:	Epp Townat:	2					×				TBA				Analysis	2	ervativ		Sampler(s):	Location:	SHE Name	C. L. N.	Project No:	
Ambient 🗆 Iced	Condition upon receipt			Manaat:						×					yl chlo anol	ride,	MTBE -	ysis +	1	List Preservative Code below:		(s):	E L	į.	Š.	ο. I	
			aadams(×					PAI	Is and	total	phenol [,]	ų.	11	below							N S >
X Refrigerated	Custody Seals		@gulfoil					×				_	-	Fec	al Coli	form			10	-		A A	281 Ea				II TATs fin. 24-l amples o
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			ill@gul		ti	 	-]	<u> </u>		CI K				To the same of the			f Chale	Gulf C	to labor cation no after 60
Di VOA Frozen Soil Jar Frozen	Present Intact Broken	*	aadams@gulfoil.com, cgill@gulfoil.com	9		æ	Group 2 PAHs - 5 µg/L	Group 1 PAHs - 0.1 µg/L	ethanol - 400 µg/L	naphthalene and vinyl chl - 5 μg/L	BTEX - 2 µg/L; TBA - 10 µg/L;	Required Minimum Levels:	* Report phenol down to MDL	State-specific reporting standards:	II*		No QC	MA DEP MCP CAM Report? Wes Wes		QA/QC Reporting Notes: * additional charges may appoly		Culobs	Chelsea State: MA	Sca - Cilillica	Gulf Chekea Terminal	Gulf Chelsea	All TATs subject to laboratory approval Min. 24-hr notification needed for rushes Samples disposed after 60 days unless otherwise instructed.
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0=01 DW=Dinking Water 7=CH3OH 8=NaHSO₄ 9=Deionized Water 10=H₃PO₄ F=Field Filtered Project Mgr: Telephone #: Report To: Andrew Adams Lab ID: X1= SO=Soil SPECTRUM ANALYTICAL, INC. Chelsea, MA 02150 281 Eastern Ave Gulf Oil LP HANIBAL TECHNOLOGY G= Grab 1=Na₂S2O₃ SL=Sludge GW=Groundwater Sample ID: Outfall 003 Outfall 003 2=HCI Andrew Adams 617.884.5980 A=Indoor/Ambient Air SW=Surface Water 3=H2SO4 3-73 4=HNO3 Date: C=Compsite Received by: SG=Soil Gas ww=Waste Water = 5=NaOH CHAIN OF CUSTODY RECORD 888 0000 Time: Invoice To: Christopher Gill P.O No.: 6=Ascorbic Acid 80 William St, Suite 400 Wellesley, MA 02481-3705 Gulf Oil LP 9 G Туре 12= WS WS Matrix Date: of VOA Vials of Amber Glass Quote/RQN: Containers # of Clear Glass Time: 1:00 # of Plastic Total Recov. (Cd, Cr, none Acid C Cu, Pb, Ni, Zn)* LC50 ** List Preservative Code below: Condition upon receipt X Ambient 🗆 Iced E-mail to: EDD format: Sampler(s): Location: Site Name: Project No: aadams@gulfoil.com, cgill@gulfoil.com ☐ Rush TAT - Date Needed: Standard TAT - 7 to 10 business days ☐ Refrigerated Custody Seals: All TATs subject to laboratory approval
Min. 24-hr notification needed for rushes
Samples disposed after 60 days unless otherwise instructed 281 Eastern Ave, Chelsen Special Handling: Gulf Chelsea Terminal ☐ Check if chlorinated ☐ DI VOA Frozen Gulf Chelsea Present CT DPH RCP Report? ☐ ASP A*
☐ NJ Reduced*
☐ Tier II* * Report metals down to MDL ✓ Standard Required Minimum Levels: State-specific reporting standards: * additional charges may appply QA/QC Reporting Notes: Cd, Pb, Ni - 0.2 ug/L **LC50 sub to GZA □ DQA* Cu - 0.5 µg/L Zn - 5 µg/L Cr - 1 µg/L Intact □ NJ Full*
□ Tier IV* ☐ ASP B* ☐ No QC Soil Jar Frozen State: Broken MA

Batch Summary

'[none]'

Aquatic Toxicity

SC32731-01 (Chelsea Creek) SC32732-01 (Outfall 003)

1705063

Microbiological Analyses

SC32732-01 (Outfall 003)

1705073

Semivolatile Organic Compounds by GCMS

1705073-BLK1 1705073-BLK2 1705073-BS1 1705073-BS2 1705073-BSD1 1705073-BSD2 SC32731-01 (Chelsea Creek)

SC32732-01 (Outfall 003)

1705079

General Chemistry Parameters

1705079-SRM1 1705079-SRM2

SC32731-01 (Chelsea Creek) SC32732-01 (Outfall 003)

<u>1705084</u>

General Chemistry Parameters

1705084-BLK1 1705084-BS1 1705084-DUP1 1705084-MS1 1705084-MSD1 1705084-SRM1 SC32731-01 (Chel

SC32731-01 (Chelsea Creek) SC32732-01 (Outfall 003)

1705104

Volatile Organic Compounds

1705104-BLK1 1705104-BS1 1705104-BSD1 SC32731-01 (Chelsea Creek) SC32732-01 (Outfall 003)

1705113

General Chemistry Parameters

1705113-BLK1 1705113-BS1 1705113-DUP1 SC32731-01 (Chelsea Creek) SC32732-01 (Outfall 003)

1705119

Total Metals by EPA 200/6000 Series Methods

SC32731-01 (Chelsea Creek) SC32732-01 (Outfall 003)

1705143

Total Metals by EPA 200 Series Methods

1705143-BLK1 1705143-BS1 1705143-DUP1 1705143-MS1 SC32731-01 (Chelsea Creek) SC32732-01 (Outfall 003)

1705144

Total Metals by EPA 200 Series Methods

1705144-BLK1 1705144-BS1 1705144-DUP1 1705144-MS1 1705144-PS1 SC32731-01 (Chelsea Creek) SC32732-01 (Outfall 003)

1705173

General Chemistry Parameters

1705173-BLK1 1705173-BS1 1705173-MS1 1705173-SRM1 SC32731-01 (Chelsea Creek) SC32732-01 (Outfall 003)

1705207

General Chemistry Parameters

1705207-BLK1 1705207-BS1 SC32731-01 (Chelsea Creek) SC32732-01 (Outfall 003)

1705216

General Chemistry Parameters

1705216-SRM1 1705216-SRM2 SC32731-01 (Chelsea Creek) SC32732-01 (Outfall 003)

1705454

General Chemistry Parameters

1705454-BLK1 1705454-BS1 1705454-DUP1 1705454-MS1 1705454-MSD1

1705454-SRM1

SC32731-01 (Chelsea Creek) SC32732-01 (Outfall 003)

380670A

Subcontracted Analyses

BX93085-BLK BX93085-LCS

SC32732-01 (Outfall 003)

S701962

Semivolatile Organic Compounds by GCMS

S701962-CAL1 S701962-CAL2 S701962-CAL3 S701962-CAL4 S701962-CAL5 S701962-CAL6 S701962-CAL7 S701962-CAL8 S701962-CAL9 S701962-CALA S701962-ICV1 S701962-ICV1 S701962-LCV1 S701962-LCV2

S702752

Semivolatile Organic Compounds by GCMS

S702752-CAL1

S701962-LCV3 S701962-TUN1

S702752-CAL2

S702752-CAL3 S702752-CAL4

S702752-CAL5

S702752-CAL6

S702752-CAL7

S702752-CAL8

S702752-CAL9

S702752-ICV1

S702752-LCV1

S702752-LCV2

S702752-TUN1

S703239

Volatile Organic Compounds

S703239-CAL1

S703239-CAL2 S703239-CAL3 S703239-CAL4 S703239-CAL5 S703239-CAL6 S703239-CAL7 S703239-CAL8 S703239-CAL9 S703239-CALA S703239-CALB S703239-ICV1 S703239-ICV1 S703239-LCV2 S703239-LCV3 S703239-LCV3

S703263

Volatile Organic Compounds

S703263-CCV1 S703263-TUN1

S703276

Semivolatile Organic Compounds by GCMS

S703276-CCV1 S703276-TUN1

S703277

Semivolatile Organic Compounds by GCMS

S703277-CCV1 S703277-TUN1